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To all men and women whose passion are algorithms, data structures, and combinatorics on strings.
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Preface

The Annual Symposium on Combinatorial Pattern Matching (CPM) has by now over 30 years of tradition and is considered to be the leading conference for the community working on Stringology. The objective of the annual CPM meetings is to provide an international forum for research in combinatorial pattern matching and related applications such as computational biology, data compression and data mining, coding, information retrieval, natural language processing, and pattern recognition.

This volume contains the papers presented at the 34th Annual Symposium on Combinatorial Pattern Matching (CPM 2023) held on June 26–28, 2023 in Marne-la-Vallée, France. The conference program includes 26 contributed papers and three invited talks, by

- Olgica Milenkovic (University of Illinois Urbana-Champaign, USA),
- Tatiana Starikovskaya (École Normale Supérieure, France), and
- Virginia Vassilevska Williams (Massachusetts Institute of Technology, USA).

For the fifth time, CPM includes the “Highlights of CPM” special session, for presenting the highlights of recent developments in combinatorial pattern matching. In this fifth edition we selected as highlight papers “Separating words and trace reconstruction”, by Zachary Chase, presented at STOC 2021, and “Approximating Dynamic Time Warping Distance Between Run-Length Encoded Strings”, by Zoe Xi and William Kuszmaul, presented at ESA 2022. The conference was preceded by a one-day student summer school taught by Karel Brinda (Inria/IRISA Rennes) and Panagiotis Charalampopoulos (Birkbeck, University of London), and organized at the École Normale Supérieure in Paris by Gabriel Bathie, Pawel Gawrychowski, Garance Gourdel and Tatiana Starikovskaya.

The contributed papers were selected out of 44 submissions, corresponding to an acceptance ratio of 59%. Each submission received at least three reviews. We thank the members of the Program Committee and all the additional external subreviewers, who are listed below, for their hard, invaluable, and collaborative effort that resulted in an excellent scientific program.

The Annual Symposium on Combinatorial Pattern Matching started in 1990, and has since then taken place every year. Previous CPM meetings were held in Paris, London (UK), Tucson, Padova, Asilomar, Helsinki, Laguna Beach, Aarhus, Piscataway, Warwick, Montreal, Jerusalem, Fukukoa, Morelia, Istanbul, Jeju Island, Barcelona, London (Ontario, Canada), Pisa, Lille, New York, Palermo, Helsinki, Bad Herrenalb, Moscow, Ischia, Tel Aviv, Warsaw, Qingdao, Pisa, Copenhagen (on-line), Wroclaw and Prague. From 1992 to the 2015 meeting, all proceedings were published in the LNCS (Lecture Notes in Computer Science) series. Since 2016, the CPM proceedings have appeared in the LIPIcs (Leibniz International Proceedings in Informatics) series, as volume 54 (CPM 2016), 78 (CPM 2017), 105 (CPM 2018), 128 (CPM 2019), 161 (CPM 2020), 191 (CPM 2021) and 223 (CPM 2022). The entire submission and review process was carried out using the EasyChair conference system.

We thank the CPM Steering Committee for their support and advice.

Laurent Bulteau and Zsuzsanna Lipták
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Trie-Compressed Adaptive Set Intersection

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Abstract

We introduce space- and time-efficient algorithms and data structures for the offline set intersection problem. We show that a sorted integer set \( S \subseteq [0..u) \) of \( n \) elements can be represented using compressed space while supporting \( k \)-way intersections in adaptive \( O(k\delta \log(u/\delta)) \) time, \( \delta \) being the alternation measure introduced by Barbay and Kenyon. Our experimental results suggest that our approaches are competitive in practice, outperforming the most efficient alternatives (Partitioned Elias-Fano indexes, Roaring Bitmaps, and Recursive Universe Partitioning (RUP)) in several scenarios, offering in general relevant space-time trade-offs.

2012 ACM Subject Classification Theory of computation → Data compression; Theory of computation → Design and analysis of algorithms; Theory of computation → Data structures and algorithms for data management; Information systems → Information retrieval query processing

Keywords and phrases Set intersection problem, Adaptive Algorithms, Compressed and compact data structures

1 Introduction

Sets are one of the most fundamental mathematical concepts related to the storage of data. Operations such as set intersections, unions, and differences are key for querying them. E.g., the use of logical AND and OR operators in web search engines translate into intersections and unions, respectively. Representing sets to support their basic operations efficiently has been a major concern since many decades ago [4]. In several applications, such as query processing in information retrieval (IR) [15] and database management systems (DBMS) [23], sets are known in advance to queries, hence data structures can be built to speed up query processing. With this motivation, in this paper we focus on the following problem.

**THE OFFLINE SET INTERSECTION PROBLEM, OSIP**

- **Input:** A family \( S = \{S_1, \ldots, S_N\} \) of \( N \) sorted integer sets over universe \([0..u), \) with \(|S_i| = n_i\).
- **Task:** To preprocess family \( S \) to efficiently support query instances of the form \( Q = \{i_1, \ldots, i_k\} \subseteq \{1..N\} \), which ask to compute \( I(Q) = \bigcap_{i \in Q} S_i \).
We assume $u = 2^k$ in this paper, for $k \geq 0$. Unless explicitly otherwise stated, we also assume $\lg x = \lfloor \lg x \rfloor$ and $\lg 0 = 0$. Typical applications of this problem include the efficient support of join operations in DBMS [23, 51], query processing using inverted indexes in IR [15, 52], and computational biology [33], among others. Building a data structure to speed up intersections, however, increases the space usage. Today, data-intensive applications encourage not only time- but also space-efficient solutions [7]. Being able to process big datasets entirely in main memory is the main motivation. Compact, succinct, and compressed data structures are important to achieve this [41]. We study here compressed data structures to efficiently support the OSIP. We assume the word RAM model of computation with word size $w = \Theta(\lg u)$. Arithmetic, logic, and bitwise operations, as well as accesses to $w$-bit memory cells, take $O(1)$ time.

The literature on this problem is vast. For the online version of the problem, where sets to be intersected are given at query time – so there is no time to preprocess them – algorithms like the ones by Baeza-Yates [9], Demaine et al. [20], and Barbay and Kenyon [12] are among the most efficient and well-known approaches. In particular, the two latter algorithms are adaptive, meaning that they are able to perform faster on “easier” query instances. The algorithm by Barbay and Kenyon runs in optimal $O(\delta \sum_{i \in Q} \lg(n_i/\delta))$ time, where $\delta$ is the so-called alternation measure that quantifies the query difficulty [12]. The algorithm by Demaine et al. [20] has running time $O(k\delta \lg(n/\delta))$, for $n = \sum_{i \in Q} n_i$, which is optimal when $\max_{i \in Q} \{\lg n_i\} = O(\min_{i \in Q} \{\lg n_i\})$ [11]. These algorithms require sets to be stored in plain form, e.g. using a sorted array or a B-tree [20], requiring $\Theta(mw)$ bits of space, for $m = \sum_{i=1}^N n_i$. This can be excessive when dealing with large databases.

For the OSIP, we have the extensive literature on inverted indexes [52, 55, 15, 46], whose main focus is on practical space-efficient set representations supporting intersections. Approaches like Optimized PForDelta [53], Roaring Bitmaps [36], SIMD-BP128 [35], and Recursive Universe Partitioning [45] shine in practical scenarios, yet without appealing theoretical guarantees of space usage and intersection computation time. Another relevant approach on these lines is Partitioned Elias-Fano (PEF) [43], able to exploit the distribution and clustering of set elements to improve space usage. Barbay and Kenyon’s algorithm can be implemented on PEF, taking $O(\delta \sum_{i \in Q} \lg(u/n_i))$ time. Regarding space usage, there is no known bound (although it performs well in practice). On a more theoretical track, Bille et al. [14] introduce a data structure that uses $O(mw)$ bits of space and supports intersections in $O(n \lg^2(w/w + k|Q|))$ time. Cohen and Porat [18] data structure also uses $O(mw)$ bits of space and allows one to compute the intersection between any two sets in $S$ in $O(\sqrt{N|Q|} + |Q|)$ time. Besides using linear space, this approach only works for pair-wise intersections (and is hard to efficiently extend to multiway intersections). Finally, Ding and Konig [21] introduce a data structure able to compute intersections in $O(n/\sqrt{w} + k|Q|)$ expected time, and uses linear $O(m)$ space. The space can be improved in practice to use about 1.88 times the space of an Elias $\gamma/\delta$ compressed inverted index [21], yet with no theoretical guarantees. Later, Gagie et al. [26] showed that wavelet trees [28] can support intersections in $O(k\delta \lg(u/\delta))$ time, using uncompressed $mw(1 + o(1))$ bits of space.

In this paper we show that $O(k\delta \lg(u/\delta))$ intersection time using compressed space is possible. In particular, (1) in Section 3 we revisit a classic (and neglected) algorithm by Trabb-Pardo [50] (former Knuth’s student) to prove that its running time is actually $O(k\delta \lg(u/\delta))$ – so it is likely the first adaptive intersection algorithm that ever existed; (2) in Section 4 we show that Trabb-Pardo’s algorithm can be implemented in compressed space, yielding an adaptive and compressed set intersection algorithm; (3) in Section 5 we show how to exploit the presence of runs of successive elements, typical in some applications [8],
to formally improve both space usage of input sets and the intersection computation time by introducing an intersection algorithm that runs in time $O(k\xi \lg (u/\xi))$, where $\xi \leq \delta$ is an adaptability measure we introduce; and (4) in Sections 6 and 7 we implement our proposals and show preliminary experimental results that indicate that our approaches are appealing not only in theory, but also in practice, outperforming the most competitive state-of-the-art approaches in some practical inverted-index datasets we use in our tests. Overall, we conclude that both theoretical guarantees and practicality can be achieved with a single approach, which is a step forward in bridging the gap between theory and practice in this important line of research.

\section{Preliminaries and Related Work}

\subsection{Operations \text{rank} and \text{select}}

The following operations on a sorted integer set $S$ are of interest:

- $\text{rank}(S, x)$: for $x \in [0..u)$, yields $\lfloor \mathsf{y} \in S \mid y \leq x \rfloor$.
- $\text{select}(S, j)$: for $1 \leq j \leq |S|$, yields $x \in S$ s.t. $\text{rank}(S, x) = j$.

A set $S$ can be alternatively described using its characteristic bit vector (cbv, for short) $C_S[0..u)$, such that $C_S[x] = 1$ if $x \in S$, $C_S[x] = 0$ otherwise. On a cbv $C_S$ we define:

- $C_S.\text{rank}_1(x)$: for $x \in [0..u)$, yields the number of 1s in $C_S[0..x]$.
- $C_S.\text{select}_1(k)$: for $1 \leq k \leq |S|$, yields the smallest position $0 \leq x < u$ s.t. $C_S.\text{rank}_1(x) = k$.

Notice that $\text{rank}(S, x) \equiv C_S.\text{rank}_1(x)$ and $\text{select}(S, j) \equiv C_S.\text{select}_1(j)$.

\subsection{Set Compression Measures}

A compression measure quantifies the amount of bits needed to encode data using a particular compression model. For an integer universe $U = [0..u)$, let $C^{(n)} \subseteq 2^U$, $n \in U$, denote the class of all sets $S \subseteq U$ such that $|S| = n$. We assume $S = \{x_1, \ldots, x_n\}$, for $0 \leq x_1 \leq \cdots \leq x_n < u$. As $|C^{(n)}| = \binom{u}{n}$, in the worst case one needs at least $B(n, u) = \lceil \lg \binom{u}{n} \rceil$ bits to encode a set $S \in C^{(n)}$. If $n \ll u$, $B(n, u) = n \lg (u/n) + n \lg e - O(\lg u)$ bits (using Stirling for $n!$). Notice $B(n, u)$ is a worst-case lower bound: some sets in $C^{(n)}$ can be encoded using less bits, as we shall see.

\subsubsection{The $\text{gap}(S)$ Compression Measure}

Let us denote $g_1 = x_1$ and, for $i = 2, \ldots, n$, $g_i = x_i - x_{i-1} - 1$. Thus, in the gap model we have $C_S[0..u) = 0^u \cdot 1^e \cdot 0^{g_1} \cdot 1^f \cdot 0^{g_2} \cdot 1^f \cdot \cdots \cdot 0^{g_n} \cdot 1^f$ (assuming wlog that $C_S$ ends with 1). Then, we define $\text{gap}(S) = \sum_{i=1}^n (\lfloor \lg g_i \rfloor + 1)$, as the amount of bits required to represent $S$ provided we encode the sequence of gaps $G = (g_1, \ldots, g_n)$, using $\lceil \lg g_i \rceil + 1$ bits per gap. Although this measure is not achievable, it exploits the variation in the gaps between consecutive set elements: the closer the elements, the smallest this measure is. It holds that $\text{gap}(S) \leq n \lg \frac{n}{\delta}$, with equality only when $g_i = \frac{u-i}{n}$ (for $i = 1, \ldots, n$). This is a measure traditionally used in applications like inverted-index compression in information retrieval [15] and databases [52].

\subsubsection{The $\text{rle}(S)$ Compression Measure}

When set elements tend to be clustered into runs of successive elements, a (usually) better way to model its cbv is $C_S[0..u) = 0^{\ell_1} \cdot 1^e \cdot 0^{\ell_2} \cdot 1^f \cdot \cdots \cdot 0^{\ell_n} \cdot 1^f$, where the sequences $Z = (z_1, \ldots, z_r)$ and $O = (\ell_1, \ldots, \ell_r)$ are the lengths of the alternating 0/1-runs in $C_S$ (assume wlog that $C_S$ begins with 0 and ends with 1). Then, $\text{rle}(S) = \sum_{i=1}^r (\lfloor \lg (z_i - 1) \rfloor + 1) + \sum_{i=1}^r (\lfloor \lg (\ell_i - 1) \rfloor + 1)$. Unfortunately, $\text{gap}(S)$ and $\text{rle}(S)$ are not comparable measures. If $n < u/2$, it holds that $\text{rle}(S) < B(n, u) + n + O(1)$ [24].
2.2.3 The \(\text{trie}(S)\) Compression Measure

Let us consider now representing a set \(S \in C(n)\) using a binary trie denoted \(\text{bintrie}(S)\), where the \(\ell = \lfloor \log u \rfloor\)-bit binary encoding of every element is added. Each internal node in \(\text{bintrie}(S)\) has two children, the left one corresponding to bit 0 and the right one to bit 1. The external nodes of \(\text{bintrie}(S)\) have no children, as usual. In our case, we distinguish two kinds of external nodes. A void external node is one whose depth is either \(d < \ell\), or alternatively \(d = \ell\) yet it represents no element in \(S\). A valid external node (or, simply, external node, or alternatively a leaf), on the other hand, is one whose depth is exactly \(\ell\) and corresponds to an element in \(S\). Thus, \(\text{bintrie}(S)\) has \(|S|\) valid external nodes, all at depth \(\ell\). For a leaf \(v\) corresponding to element \(x_i \in S\), the root-to-\(v\) path is hence labeled with the binary encoding of \(x_i\). This approach has been used for representing sets since at least the late 70s by Trabb-Pardo [50]. Consider the example sets \(S_1 = \{1, 3, 7, 8, 9, 11, 12\}\) and \(S_2 = \{2, 5, 7, 12, 15\}\) over universe \([0, 16]\), that we shall use as running examples. Figure 1 shows the corresponding tries \(\text{bintrie}(S_1)\) and \(\text{bintrie}(S_2)\), with external nodes shown as squares and void external nodes with dotted lines. Interestingly, the following compression measure can be derived from this representation [29]. Given two bit strings \(x\) and \(y\) of \(\ell\) bits each, let \(x \odot y\) denote the bit string obtained after removing the longest common prefix among \(x\) and \(y\) from \(x\). For instance, for \(x = 0110100\) and \(y = 0111011\), we have \(x \odot y = 0100\). The prefix omission method by Klein and Shapira [31] represents a sorted set \(S\) as a binary sequence \(T = (x_1; x_2 \odot x_1; \ldots ; x_n \odot x_{n-1})\). If we denote \(|x_i \odot x_{i-1}|\) the length of bit string \(x_i \odot x_{i-1}\), then the whole sequence uses

\[
\text{trie}(S) = |x_1| + \sum_{i=2}^{n} |x_i \odot x_{i-1}|.
\]

It turns out that \(\text{trie}(S)\) is the number of edges in \(\text{bintrie}(S)\) [29]. Notice that \(\text{trie}(S)\) decreases as longer trie paths are shared among set elements: consider two integers \(x\) and \(y\), the trie represents their longest common prefix just once (then saving space), and then represents both \(x \odot y\) and \(y \odot x\). Extreme cases are as follows: (1) All set elements form a single run of consecutive elements, which maximizes the number of trie edges shared among set elements, hence minimizing the space usage; and (2) The \(n\) elements are uniformly distributed within \([0, u]\) (i.e., the gap between successive elements is \(g_i = u/n\)), which minimizes the number of trie edges shared among elements, and hence maximizes space usage. Notice this is similar to the case that maximizes the \(\text{gap}(S)\) measure.

**Definition 1.** We say that a node \(v\) in \(\text{bintrie}(S)\) covers all leaves that descend from it. In such a case, we call \(v\) a cover node of the corresponding leaves.
The following lemma summarizes several results that shall be important for our work:

**Lemma 2** ([26], Lemmas 1–5). For bintrie($S$), the following results hold:
1. Any contiguous range of $L$ leaves in bintrie($S$) is covered by $O(\lg L)$ nodes.
2. Any set of $r$ nodes in bintrie($S$) has $O(r \lg \frac{n}{r})$ ancestors.
3. Any set of $r$ nodes in bintrie($S$) minimally covering a contiguous range of leaves in the trie has $O(r + \lg u)$ ancestors.
4. Any set of $r$ nodes in bintrie($S$) minimally covering $L$ contiguous leaves has $O(\lg u + r \lg \frac{u}{r})$ ancestors.

**Definition 3.** Given a set $S = \{x_1, \ldots, x_n\} \subseteq [0..u)$, let $S + a$, for $a \in [0..u)$, denote a shifted version of $S$: $S + a = \{(x_1 + a) \mod u, (x_2 + a) \mod u, \ldots, (x_n + a) \mod u\}$.

The following result is relevant for our proposal:

**Lemma 4** ([29], Section 2). Given a set $S \subseteq [0..u)$ of $n$ elements, it holds that:
1. $\text{trie}(S) \leq \min\{2 \text{gap}(S), n \lg (u/n) + 2n - 2\}$.
2. $\exists a \in [0..u)$, such that $\text{trie}(S + a) \leq \text{gap}(S) + 2n - 2$.
3. $\text{trie}(S + a) \leq \text{gap}(S) + 2n - 2$ on average over all values of $a \in [0..u)$.

### 2.3 Adaptive Set Intersection Algorithms

An adaptive algorithm is one whose running time is a function not only of the instance size (as usual), but also of a difficulty measure of the instance. In this way, “easy” instances are solved faster than “difficult” ones, allowing for a more refined analysis than typical worst-case approaches. For the set intersection problem, algorithms by Demaine, López-Ortiz, and Munro [20] and by Barbay and Kenyon [12] are the most important adaptive approaches. To analyze adaptive intersection algorithms, Demaine et al. and Barbay and Kenyon agree in that any algorithm that computes $I(Q)$ must show a certificate [12] or proof [20] to prove that the intersection is correct. That is, that any element in $I(Q)$ belongs to the $k$ sets $S_1, \ldots, S_k$, and no element in the intersection has been left out of the result. Then, the analysis determines the size of a certificate (or proof) and the time it takes to compute them. In particular, Barbay and Kenyon [12] partition certificates are defined as follows.

**Definition 5.** Given a query $Q = \{i_1, \ldots, i_k\} \subseteq [1..N]$, a partition certificate is a partition of the universe $[0..u)$ into a set of intervals $P_\text{br}(Q) = \{I_1, I_2, \ldots, I_p\}$, such that:
1. $\forall x \in I(Q), [x..x] \in P_\text{br}(Q)$;
2. $\forall x \notin I(Q), \exists I_j \in P_\text{br}(Q), x \in I_j \land \exists q \in Q, S_q \cap I_j = \emptyset$.

For a given query $Q$, several valid partition certificates could be given. However, we are interested in the smallest partition certificate of $Q$, as it takes the least time to be computed.

**Definition 6.** For a given query instance $Q = \{i_1, \ldots, i_k\} \subseteq [1..N]$, let $\delta$ denote the size of the smallest partition certificate of $Q$.

Measure $\delta$ is known as the alternation of the query instance [12], measuring its difficulty. Notice $|I(Q)| \leq \delta$ holds. Figure 2 shows the smallest partition certificate (of size $\delta = 8$) for sets $S_1$ and $S_2$ of our running example. Barbay and Kenyon [11, 12] proved a lower bound of $\Omega(\delta \sum_{i \in Q} \lg (n_i/\delta))$ comparisons for the set intersection problem. They also gave and optimal intersection algorithm running in $O(\delta \sum_{i \in Q} \lg (n_i/\delta))$ time.
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$S_1$: 1 3 5 7 8 9 10 11 12
$S_2$: 2 5 7 12 15

**Figure 2** Vertical lines show the smallest partition certificate $P = \{[0..1], [2..2], [3..4], [5..6], [7..7], [8..11], [12..12], [13..15]\}$ of size $\delta = 8$ of the universe $[0..16)$ for the intersection of sets $S_1 = \{1, 3, 7, 8, 9, 10, 11, 12\}$ and $S_2 = \{2, 5, 7, 12, 15\}$.

3 Trie Intersection Certificates: A Revisit to Trabb-Pardo Algorithm

In this section we revisit an old divide-and-conquer intersection algorithm by Trabb-Pardo [50], not only to review it but also to prove an adaptive bound on its running time. Algorithm 1 shows the pseudocode. Given a query instance $Q = \{i_1, \ldots, i_k\} \subseteq [1..N]$, the algorithm must be invoked as $\text{TP-Intersection}(S_{i_1}, \ldots, S_{i_k}, [0..u])$. The main idea is to divide the universe into two halves, to then split each set according to this universe division. This differs from, e.g., Baeza-Yates’s algorithm [9, 10], which splits according to the median of one of the sets. The Divide steps (lines 10 and 11) can be implemented using binary search. At the first level of recursion, the most-significant bit of every element in sets $S_{i,l}$ is 0, as they belong to the left half of the universe. Similarly, for $S_{i,r}$, the most-significant bit is 1 as all elements belong to the right half. At each node of the recursion tree, the current universe is divided into two halves, to then recurse on the sets split accordingly.

As sets are known in advance to queries and set splits carried out by Algorithm 1 depend just on the universe, the Divide step of Algorithm 1 can be implemented efficiently by using a suitable set representation that not only stores the set values, but also precomputes the set splits carried out recursively by the algorithm. Trabb-Pardo proposes to represent each $S_i \in S$ using $\text{bintrie}(S_i)$, mimicking the way set $S_i$ is recursively split by Algorithm 1. The left child of the root represents all elements whose most-significant bit is 0, i.e., elements in set $S_{i,l}$ of Algorithm 1 (line 10) in the first level of recursion; similarly for $S_{i,r}$, containing...
all elements in $S_i$ whose most-significant bit is 1. To simulate the recursive execution of Algorithm 1 on the binary tries, one must carry out a DFS traversal in synchronization on all tries involved in the query, following the same path in all of them and stopping (and backtracking if needed) as soon as we reach a dead end in one of the tries (which correspond to dotted lines in Figure 1), or we reach a leaf node in all the tries (in which case we have found an element belonging to the intersection). In this way, (1) we stop as soon as we detect a universe interval that does not have any element in the intersection, and (2) we find the relevant elements when we arrive at the leaves.

To analyze Algorithm 1, we introduce the concept of trie intersection certificates, denoted $\text{cert}(Q)$, as an alternative to existing certificates [20, 12]. Figure 3 shows a possible $\text{cert}(Q)$ for the intersection of $S_1$ and $S_2$ from Figure 1. Let $\text{path}(v)$ denote the binary string labelling the root-to-$v$ path, and $\text{depth}(v) = |\text{path}(v)|$. For a query $Q$, a binary trie $\text{cert}(Q)$ is a trie partition certificate if: (1) for any internal node $v$ of $\text{cert}(Q)$ such that $\text{path}(v) = b$, there exists an internal node $v_i$ with $\text{path}(v_i) = b$ in every $\text{bintrie}(S_i)$, $i \in Q$; (2) for any void external node $v$ with $\text{depth}(v) = d \leq \ell = \lceil \log u \rceil$ and $\text{path}(v) = b$, there exists at least a set $S_i$ ($i \in Q$) such that there is no node $v_j$ with $\text{path}(v_j) = b$ in $\text{bintrie}(S_i)$. So, the universe interval $[\text{dec}(b \cdot 0^\ell d) .. \text{dec}(b \cdot 1^\ell d)]$ has no element in the intersection, where $\text{dec}(x)$ denotes the decimal representation of a binary string $x$. We call them fail nodes, shown as “×” in Figure 3; and (3) for any valid external node $v$ with $\text{depth}(v) = \ell$ corresponding to $\text{path}(v) = b$, there exists a valid external node $v_i$ with $\text{path}(v_i) = b$ in every $\text{bintrie}(S_i)$, $i \in Q$. We call them success nodes as they correspond to elements in $I(Q)$.

Notice that $\text{cert}(Q)$ is the trie obtained by intersecting $\text{bintrie}(S_i)$, for all $i \in Q$, and that it is the smallest trie that allows us to prove the correctness of the intersection. For instance, Figure 3 shows $\text{cert}(Q)$ for a given query. Interestingly, the recursion tree of Algorithm 1 is exactly $\text{cert}(Q)$, as the algorithm stops as soon as one arrives at a fail node. The external nodes of $\text{cert}(Q)$ cover the universe $[0..u]$ with intervals, similar to Barbay and Kenyon partition certificates, as stated by the following definition.

**Definition 7.** Given a query $Q = \{i_1, \ldots, i_k\} \subseteq [1..N]$, its trie partition certificate is a partition of the universe $[0..u]$ into a set of intervals, we say that $\text{cert}(Q)$ induces the following partition of $[0..u]$ into a set of intervals that we call trie partition certificate:

\[
P_{\text{PC}}(Q) = \bigcup_{l \in \mathcal{E}(\text{cert}(Q))} \left\{ [\text{dec}(\text{path}(l)) \cdot 0^{\ell - \text{depth}(l)}) .. \text{dec}(\text{path}(l) \cdot 1^{\ell - \text{depth}(l)})] \right\},
\]

where $\mathcal{E}(\text{cert}(Q))$ denotes the set of external nodes of $\text{cert}(Q)$.

For instance, the trie certificate of Figure 3 induces the following trie partition certificate of the universe $[0..16]$: \{[0..1], [2..2], [3..3], [4..5], [6..6], [7..7], [8..11], [12..12], [13..13], [14..15]\}.
Since \( \text{cert}(Q) \) is the recursion tree of Algorithm 1, its running time is \( O(k|\text{cert}(Q)|) \). As in the worst-case one must traverse completely all tries \( \text{bintrie}(S_i), i \in Q \), we have:

\[
k|\text{cert}(Q)| \leq \sum_{i \in Q} \text{trie}(S_i) \leq \sum_{i \in Q} n_i \frac{u}{n_i} + 2n_i - 2,
\]

where the last bound is from Lemma 4 (1). Next we prove an adaptive bound for \( k|\text{cert}(Q)| \).

**Theorem 8.** Given a query instance \( Q = \{i_1, \ldots, i_k\} \subseteq [1..N] \) with alternation measure \( \delta \) and over sets with universe \([0..u]\), algorithm TP-Intersection computes \( I(Q) = \cap_{i \in Q} S_i \) in time \( O(k\delta \lg(u/\delta)) \).

**Proof.** Consider a smallest partition certificate \( \mathcal{P}_{br}(Q) = \{I_1, \ldots, I_\delta\} \) of universe \([0..u]\), such that \(|I_i| = L_i\) for \( i = 1, \ldots, \delta \). Let us think now of the worst-case smallest \( \text{cert}(Q) \) we could have, by covering the \( \delta \) intervals in \( \mathcal{P}_{br}(Q) \) with as many external nodes of \( \text{cert}(Q) \) as possible. For any \( I_j \in \mathcal{P}_{br}(Q) \) formed by elements not in \( I \), there exists a set of external fail nodes in \( \text{cert}(Q) \) that cover \( I_j \). This is because when traversing the tries \( \text{bintrie}(S_i) \) in coordination, \( i \in Q \), the algorithm stops as long as one gets into one of the cover nodes of \( I_j \), since it does not belong to at least one of the tries. According to Lemma 2 (1), a contiguous range of \( L \) leaves (corresponding to the values in \( I_j \)) can be covered with up to \( O(\lg L) \) nodes. Thus, in the worst-case, \( \text{cert}(Q) \) has \( O(\sum_{i=1}^\delta \lg L_i) \) external nodes that overall cover \([0..u]\). Now, recall that the external nodes of \( \text{cert}(Q) \) cover the contiguous range of leaves corresponding to \([0..u]\). Hence, according to Lemma 2 (3), these external nodes have \( O(\sum_{i=1}^\delta \lg L_i + \lg u) \) ancestors, so overall \( \text{cert}(Q) \) has \( O(\sum_{i=1}^\delta \lg L_i + \lg u) \) nodes. The sum is maximized when \( L_i = u/\delta \), for all \( 1 \leq i \leq \delta \), hence \( \text{cert}(Q) \) has \( O(\delta \lg(u/\delta)) \) nodes. The result follows from the fact that for each node in \( \text{cert}(Q) \) the algorithm runs in time \( O(k) \). ▶

## 4 Compressed Intersectable Sets

We devise next a space-efficient representation of \( \text{bintrie}(S) \), for a set \( S = \{x_1, \ldots, x_n\} \subseteq [0..u] \) of \( n \) elements such that \( 0 \leq x_1 < \cdots < x_n < u \). This representation will also allow for efficient intersections, supporting Trabb-Pardo’s [50] algorithm.

We represent \( \text{bintrie}(S) \) level-wise [30]. Let \( B_1[1..2l_1], \ldots, B_\ell[1..2l_\ell] \) be bit vectors such that \( B_i \) represents the \( l_i \) nodes at level \( i \) of \( \text{bintrie}(S) \) \((1 \leq i \leq \ell)\), from left to right. Each node is encoded using 2 bits, indicating the presence (using bit 1) or absence (bit 0) of the left and right children, respectively. In this way, the feasible codewords for trie nodes are \( 01, 10, \) and \( 11 \), whereas \( 00 \) is not a valid codeword. The codewords of all nodes at level \( i \geq 1 \) in the trie are concatenated from left to right to form \( B_i \). The \( j \)-th node at level \( i \) (from left to right) is stored at positions \( 2j-1 \) and \( 2j \). We say that \( 2j-1 \) is the position of such node in \( B_i \).

Let \( p \) be the position in \( B_i \) corresponding to a node \( v \) at level \( i \) of \( \text{bintrie}(S) \). As the nodes are stored level-wise and from left to right, the number of \( 1s \) before position \( p \) in \( B_i \) equals the number of nodes in level \( i+1 \) that are before the child(ren) of node \( v \). So, \( 2B_i.\text{rank}_1(p-1) + 1 \) yields the position of \( B_{i+1} \), where the first child of node \( v \) is. Figure 4 illustrates our representation.

The total number of \( 1s \) in the bit vectors of our representation equals the number of edges in the trie. That is, there are \( \text{trie}(S) \) \( 1s \). Besides, the trie has \( \text{trie}(S) + 1 \) internal nodes and leaves: \( n \) of them are leaves, so \( \text{trie}(S) - n + 1 \) are internal. In our representation we only need to represent the internal trie nodes. As we encode each node using 2 bits, the total space usage for \( B_1, \ldots, B_\ell \) is \( 2(\text{trie}(S) - n + 1) \) bits. On top of them we use Clark’s data structure [16] to support \( \text{rank} \) in \( O(1) \) time, adding \( o(\text{trie}(S)) \) extra bits overall.
Given a query $Q = \{i_1, \ldots, i_k\} \subseteq [1..N]$, we traverse $\text{bintrie}(S_{i_1}), \ldots, \text{bintrie}(S_{i_k})$ using a recursive DFS traversal as in Algorithm 1. Besides the query itself, our algorithm receives: (1) an integer value, $\text{level}$, indicating the current recursion level, and (2) integer values $r_1, \ldots, r_k$, indicating the current nodes in each trie, represented as the positions of these nodes within $\text{B}_{\text{level}}$. Algorithm 2 shows the pseudo-code of our adaptive and compressed algorithm to compute the compact representation for $\text{bintrie}(I(Q))$ (denoted $T_I$ in the pseudocode). The algorithm uses a binary variable $s$, initialized with 11, which stores the bitwise-and of all current node codewords (line 4). So, $s = 00$ means that recursion must stop, $s = 10$ indicates to go down only to the left, $s = 01$ just to the right, and $s = 11$ to both children.

Lines 9–13 carry out the needed computation to go down to the left child. In particular, we compute the positions of the left-subtrie roots using $\text{rank}_4$ operation. Then, in line 13 we recursively go down to the left. The result of that recursion in stored in variable $lChild$, indicating with a 1 that the left recursion yielded a non-empty intersection, 0 otherwise. A similar procedure is carried out for the right child in lines 14–21. Line 17 determines whether we have already computed the $\text{rank}_4$s corresponding to the left child. If that is not the case, we compute them in line 18. In this way, we compute only one $\text{rank}_4$ operation per traversed node in the tries, which is important in practice. Just as for the left child, we store the result of the right-child recursion in variable $rChild$ in line 21. Finally, in line 22 we determine whether the left and right recursions yielded an empty intersection or not. If both $lChild = rChild = 0$, the intersection was empty on both children, so we return 0. Otherwise, we append $lChild$ and $rChild$ to $T_I$,$\text{B}_{\text{level}}$, as that is the codeword of the corresponding node in $T_I$. Note how we actually generate the output trie $T_I$ in postorder, after we visited both children of the current nodes, despite the input is traversed in preorder. Thus, we write the output in time proportional to its size. Although the total running time is still proportional to $|\text{cert}(Q)|$, this can save important time in practice.

Besides computing $T(I(Q)$, a distinctive feature of our algorithm is that it also allows one to obtain the sequence $(\text{rank}(S_{i_1}, x), \ldots, \text{rank}(S_{i_k}, x))$, for all $x \in I(Q)$, for free (in asymptotic terms). The idea is to compute $(\text{bintrie}(S_{i_1}), B_T, \text{rank}_4(r_1), \ldots, \text{bintrie}(S_{i_k}), B_T, \text{rank}_4(r_k))$ every time the recursion reaches level $\ell$ (i.e., just before the return of line 7 in Algorithm 2). Outputting this information is important for several applications, such as cases where set elements have satellite data associated to them. For an element $x_j \in S_i$, the associated data $d_j$ is stored in an auxiliary array $D_i[1..n]$ such that $D[\text{rank}(S_i, x_j)] = d_j$. Typical applications are inverted indexes in IR (where ranking information, such as frequencies, is associated to inverted list elements), and the Leapfrog Triejoin algorithm [51] (where at each step we must compute the intersection of sets, and for each element in the intersection we must go down following a pointer associated to it).
We have proved the following theorem:

▶ **Theorem 9.** Let $S = \{S_1, \ldots, S_N\}$ be a family of $N$ integer sets, each of size $|S_i| = n_i$ and universe $[0..u)$. There exists a data structure able to represent each set $S_i$ using $2(\text{trie}(S_i) - n_i + 1) + o(\text{trie}(S_i))$ bits, such that given a query $Q = \{i_1, \ldots, i_k\} \subseteq [1..N]$, the intersection $I(Q) = \cap_{i \in Q} S_i$ can be computed in $O(k\delta \log(u/\delta))$ time, where $\delta$ is the alternation measure of $Q$. Besides, for every $x \in I(Q)$, the data structure also allows one to obtain the sequence $(\text{rank}(S_{i_1}, x), \ldots, \text{rank}(S_{i_k}, x))$ asymptotically for free.

### 5 Compressing Runs of Elements

Next, we exploit the presence of runs of successive elements in the input sets to reduce both the space usage of the binary trie representation, as well as intersection time. Runs tend to be captured by full subtrees in the corresponding binary tries. See, e.g., the full subtree whose leaves correspond to elements 8, 9, 10, 11 in the binary trie of Figure 5. Let $v$ be a bintrie($S$) node whose subtree is full. Let $\text{depth}(v) = d$. If $b = \text{path}(v)$, the $2^{d-d}$ leaves
covered by \( v \) correspond to the integer interval \([\text{dec}(b \cdot 0^{\ell_i-1})..\text{dec}(b \cdot 1^{\ell_i-1})]\). So, the subtree of \( v \) can be removed, keeping just \( v \), saving space and still being able to recover the removed elements.

\[
\sum_{i=1}^{r} (2\ell_i - 4 \log \ell_i).
\]

This immediately implies \( r_{\text{Trie}}(S) \leq \text{trie}(S) \leq 2 \cdot \text{gap}(S) \), yet we can prove tighter bounds. Assume a set \( S \) with \( r \) runs of \( \ell_1, \ldots, \ell_r \) successive elements each, respectively. The \( \ell_i \) elements of a given run correspond to \( \ell_i \) contiguous leaves in bintrie(\( S \)), which, according to Lemma 2 (item 1), are covered by at most \( 2 \cdot \lceil \log (\ell_i/2) \rceil \) nodes. This is a pessimistic case that removes the least edges, so we analyze it. Among the cover nodes, there are 2 whose subtrees have 0 edges, 2 whose subtrees have 2 edges, 2 whose subtrees have 6 edges, and so on. In general, for each \( i = 1, \ldots, \lceil \log (\ell_i/2) \rceil \), there are 2 cover nodes whose subtrees have \( 2^i - 2 \) edges. If we remove them all, the total number of edges removed is

\[
2 \sum_{i=1}^{r} \lceil \log (\ell_i/2) \rceil (2^i - 2) \leq 2\ell_i - 4 \log \ell_i.
\]

This removes the least edges belonging to full subtrees, so we can bound

\[
r_{\text{Trie}}(S) \leq \text{trie}(S) - \sum_{i=1}^{r} (2\ell_i - 4 \log \ell_i).
\]  

We can also prove the following bounds.

**Definition 10.** Let \( S \subseteq [0..u) \) be a set of \( n \) elements. We define \( r_{\text{Trie}}(S) \) as the number of edges in bintrie(\( S \)) after removing the maximal full subtrees.

**Lemma 11.** Given a set \( S \subseteq [0..u) \) of \( n \) elements, it holds that

1. \( r_{\text{Trie}}(S) \leq 2 \cdot \min \{ r_{\text{le}}(S) + \sum_{i=1}^{r} \log \ell_i, \text{gap}(S) \} \).
2. \( \exists a \in [0..u), \) such that \( r_{\text{Trie}}(S + a) \leq \min \{ r_{\text{le}}(S) - \sum_{i=1}^{r} \ell_i + 3 \sum_{i=1}^{r} \log \ell_i, \text{gap}(S) \} + 2n - 2 \).
3. \( r_{\text{Trie}}(S + a) \leq \min \{ r_{\text{le}}(S) - \sum_{i=1}^{r} \ell_i + 3 \sum_{i=1}^{r} \log \ell_i, \text{gap}(S) \} + 2n - 2 \) on average, assuming \( a \in [0..u) \) is chosen uniformly at random.

**Proof.** Since \( S \) has \( r \) runs of \( \ell_1, \ldots, \ell_r \) elements, we can rewrite \( \text{gap}(S) = \sum_{i=1}^{r} (\lceil \log (\ell_i - 1) \rceil + 1) + \sum_{i=1}^{r} (\ell_i - 1) \). As \( r_{\text{Trie}}(S) \leq \text{trie}(S) \leq 2 \cdot \text{gap}(S) \), and \( r_{\text{Trie}}(S) \leq \text{trie}(S) - \sum_{i=1}^{r} (2\ell_i - 4 \log \ell_i) \) (Equation 1), it holds that

**Figure 5** Left side, the binary trie representing set \{1, 3, 7, 8, 9, 10, 11, 12\}. Notice that the subtree whose leaves correspond to elements 8, 9, 10, 11 is a full subtree. Right side, our compact representation removing full subtrees and encoding their roots with 00.
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\[ r_{\text{Trie}}(S) \leq \text{trie}(S) - \sum_{i=1}^{r} (2\ell_i - 4 \lg \ell_i) \]

\[ \leq 2\sum_{i=1}^{r} (\lceil \lg (z_i - 1) \rceil + 1) + \sum_{i=1}^{r} (\ell_i - 1)) - \sum_{i=1}^{r} (2\ell_i - 4 \lg \ell_i) \]

\[ = 2\sum_{i=1}^{r} (\lceil \lg (z_i - 1) \rceil + 1) + 4 \sum_{i=1}^{r} \lg \ell_i = 2(r_{\text{Trie}}(S) + \sum_{i=1}^{r} \lg \ell_i), \]

proving item 1. Items 2 and 3 can be proved similarly from items 2 and 3 of Lemma 4. ▶

In our compact representation, we encode a full-subtree cover node using 00. Recall that 00 is an invalid codeword, so we use it as a special mark. See Figure 5 for an illustration.

Given a query \( Q = \{i_1, \ldots, i_k\} \subseteq [1..N] \), the procedure to compute \( I(Q) \) is similar to that of Algorithm 2. The only difference is that if in a given trie \( \text{bintrie}(S_i) \) we arrive at a node encoded 00, every possible set element in the subtrie of the node belongs to \( S_i \). In other words, the intersection within the current subtries is independent of \( S_i \), so we can safely temporarily exclude \( \text{bintrie}(S_i) \) from the intersection and continue intersecting the remaining tries. To implement this idea, we keep boolean flags \( f_1, \ldots, f_k \) such that \( f_j \) corresponds to \( \text{bintrie}(S_i) \). The idea is that at each point during the synchronized DFS traversal, only tries whose flag is true participate in the intersection. Initially, we set \( f_1 \leftarrow \text{false} \), for \( 1 \leq i \leq k \). If, during the intersection process, we arrive at a node encoded 00 in \( \text{bintrie}(S_i) \), we set \( f_i \leftarrow \text{true} \). When the recursion at a node encoded 00 in \( \text{bintrie}(S_i) \) finishes, we set \( f_i \leftarrow \text{false} \) again. If, at a given point, all tries have been temporarily excluded but one, let us say \( \text{bintrie}(S_j) \), we only need to traverse the current subtree in \( S_j \), copying it verbatim to the output. If this subtree contains nodes encoded 00, they will appear in the output. This way, the maximal runs of successive elements in the output will be covered by nodes encoded 00. This fact is key for the adaptive running time of our algorithm, as we shall see below.

We analyze our algorithm introducing the following variant of partition certificates.

▶ Definition 12. Given a query instance \( Q = \{i_1, \ldots, i_k\} \subseteq [1..N] \), a run-partition certificate for it is a partition of the universe \([0,u]\) into a set of intervals \( \mathcal{P}_{\text{rk}}(Q) = \{I_1, I_2, \ldots, I_p\} \), such that the following conditions hold:

1. \( \forall x \in I(Q), \exists I_j \in \mathcal{P}_{\text{rk}}(Q), \text{ such that } x \in I_j \land I(Q) \cap I_j = I_j; \)
2. \( \forall x \not\in I(Q), \exists I_j \in \mathcal{P}_{\text{rk}}(Q), \text{ such that } x \in I_j \land \exists q \in Q, S_q \cap I_j = \emptyset. \)

Let \( \xi \) denote the size of the smallest run-partition certificate \( \mathcal{P}_{\text{rk}}(Q) \) of \( Q \). We call \( \xi \) the run alternation measure.

Item 2 is the same as for Barbay and Kenyon’s partition certificates, corresponding to intervals of elements not in \( I(Q) \). Item 1, on the other hand, corresponds to elements in \( I(Q) \) which, unlike Barbay and Kenyon certificates, are not necessarily covered by singletons: our definition allows one to cover a run of successive elements in \( I(Q) \) using a single interval. Clearly, \( \xi \leq \delta \) holds. Besides, although \( |I(Q)| \leq \delta \) holds, in our case there can be query instances such that \( \xi < |I(Q)| \). Figure 6 illustrates our definition for an intersection of 4 sets on the universe \([0,.15] \). Notice that \( \xi = 5 \), whereas \( |I(Q)| = 6 \) and \( \delta = 9 \).

We must also introduce a fourth type of node to our trie certificate definition of Section 3. If for an internal node \( v \) of \( \text{cert}(Q) \) with \( \text{path}(v) = b \), it holds that there is a node \( v_i \) with \( \text{path}(v_i) = b \) in every \( \text{bintrie}(S_i) \), \( i \in Q \), and the subtrees of all \( 0 \)’s is full, then \( v \) is called an internal success node. It is important to note that every interval \( I_j \) from item 1 of Definition 12 is covered only by internal success nodes. Also, internal success nodes only cover intervals from item 1 of Definition 12.
the binary trie representation, and (2) the plain array representation. In our experiments we will use the latter, to be fair: all testes alternatives produce their outputs in plain form. We implemented bit vectors Algorithm 2 on our compact trie data structures, following the descriptions from Sections 4 and 5 very closely. We implemented, however, two alternatives for representing the output: (1) the binary trie representation, and (2) the plain array representation. In our experiments we will use the latter, to be fair: all testes alternatives produce their outputs in plain form.

We also implemented a simple multithreaded version of our algorithm. Let \( t \) denote the number of available threads. Then, we define \( c = \lfloor \log t \rfloor \). Our algorithm proceeds as in Algorithm 2, generating a binary trie of height \( c \) (that we will call top trie), with at most \( t \) leaves. Then, we execute Algorithm 2 again, this time in parallel, with each thread

\[
\begin{array}{c|cccccccc}
S_{i1} & 7 & 8 & 9 & 10 & 11 & 12 & 13 & 14 & 15 \\
S_{i2} & 5 & 6 & 7 & 8 & 9 & 10 & 11 & 12 & 13 & 14 \\
S_{i3} & 4 & 5 & 6 & 7 & 8 & 9 & 11 & 12 & 13 & 14 \\
S_{i4} & 8 & 9 & 10 & 11 & 12 & 13 & 14 & 15 \\
\end{array}
\]

Figure 6 A query instance \( Q = \{S_{i1}, S_{i2}, S_{i3}, S_{i4}\} \) and its smallest run-partition certificate \( \mathcal{P}_{\mathcal{C}}(Q) = \{[0..7], [8..9], [10..10], [11..14], [15..15]\} \) of size \( \xi = 5 \).

Our main result is stated in the following theorem:

> Theorem 13. Let \( S = \{S_1, \ldots, S_N\} \) be a family of \( N \) integer sets, each of size \( |S_i| = n_i \) and universe \([0..u]\). There exists a data structure able to represent each set \( S_i \) using \( 2r\text{Trie}(S_i)(1 + \mathcal{O}(\text{Trie}(S_i))) \) bits, such that given a query \( Q = \{i_1, \ldots, i_k\} \subseteq [1..N] \), the intersection \( I(Q) = \cap_{i \in Q} S_i \) can be computed in \( \mathcal{O}(k\xi \log(u/\xi)) \) time, where \( \xi \) is the run alternation measure of \( Q \).

Proof. Consider the smallest run-partition certificate \( \mathcal{P}_{\mathcal{C}}(Q) = \{I_1, \ldots, I_{\xi}\} \) of universe \([0..u]\), such that \( |I_i| = L_i \) for \( i = 1, \ldots, \xi \). Let us cover these \( \xi \) intervals with as many nodes of the smallest \( \text{cert}(Q) \) as possible. As we already saw in the proof of Theorem 8, all intervals \( I_i \) such that \( I_i \cap I(Q) = \emptyset \) are covered by at most \( \mathcal{O}(\log L_i) \) nodes in \( \text{cert}(Q) \). We now prove the same for intervals \( I_i \subseteq I(Q) \), which are covered by internal success nodes of \( \text{cert}(Q) \). The only thing to note is that our algorithm stops as soon as it arrives to an internal success node. As there can be \( \mathcal{O}(\log L_i) \) such cover nodes, universe \([0..u]\) can be covered by \( \mathcal{O}(\log u + \sum_{i=1}^{\xi} \log L_i) = \mathcal{O}(\log u + \sum_{i=1}^{\xi} \log(u/\xi)) \) nodes, hence \( \text{cert}(Q) \) has \( \mathcal{O}(\xi \log(u/\xi)) \) nodes overall. The result follows from the fact that at each node the algorithm runs in time \( \mathcal{O}(k) \).

6 Implementation

We implemented bit vectors \( B_1, \ldots, B_\ell \) in plain form using class \texttt{bit\_vector<>} from the sdsl library [27]. We support \texttt{rank} on them using different data structures to obtain the following schemes. \texttt{trie v, rTrie v}: the variants defined in Section 4 and 5, respectively, using \texttt{rank\_support\_v} for \texttt{rank1}. It uses \( \sim25\% \) extra space on top of the bit vector, supporting \texttt{rank1} in \( \mathcal{O}(1) \) time. \texttt{trie v5, rTrie v5}: use \texttt{rank\_support\_v5}, requiring \( \sim6.25\% \) extra space on top of the bit vectors, supporting \texttt{rank1} in \( \mathcal{O}(1) \) time. This alternative is smaller, yet slower in practice. \texttt{trie IL, rTrie IL}: use \texttt{rank\_support\_il}, aiming at reducing the number of cache misses to compute \texttt{rank1}. We use block size 512, requiring \( \sim12.5\% \) extra space on top of the bit vectors, while supporting \texttt{rank1} in \( \mathcal{O}(1) \) time.

Most state-of-the-art alternatives we compare with do not support operation \texttt{rank}(\( S, x \)). So, to be fair, we do not store any \texttt{rank1} data structure for the last-level bit vector \( B_\ell \). Recall that \texttt{rank}(\( S, x \)) is equivalent to a \texttt{rank1} on the corresponding position of \( B_\ell \). We implemented Algorithm 2 on our compact trie data structures, following the descriptions from Sections 4 and 5 very closely. We implemented, however, two alternatives for representing the output: (1) the binary trie representation, and (2) the plain array representation. In our experiments we will use the latter, to be fair: all testes alternatives produce their outputs in plain form.

We also implemented a simple multithreaded version of our algorithm. Let \( t \) denote the number of available threads. Then, we define \( c = \lfloor \log t \rfloor \). Our algorithm proceeds as in Algorithm 2, generating a binary trie of height \( c \) (that we will call \texttt{top trie}), with at most \( t \) leaves. Then, we execute Algorithm 2 again, this time in parallel, with each thread
starting from a different leaf of the top trie. Each thread generates its own output in parallel, using our compact trie representation. Once all threads finish, we concatenate these tries to generate the final output. We just need to count, in parallel, how many nodes there are in each level of the trie. Then, we allocate a bit vector of the appropriate size for each level, where each thread will write its own part of the output in parallel. This simple approach does not guarantee load balancing among threads, however it works relatively well in practice.

Our source code and instructions to replicate our experiments are available at https://github.com/jpcastillog/compressed-binary-tries.

7 Experimental Results

We experimentally evaluate our approaches on a server with an i7 10700k CPU, 8 cores and 16 threads at 4.70 GHz, 32 GB of RAM (DDR4-3.6GHz) running in dual channel, and Ubuntu 20.04 LTS OS. Our implementation is developed in C++, compiled with g++ 9.3.0 and optimization flags -O3 and -march=native.

In our tests, we used families of sets corresponding to inverted indexes of three standard document collections: Gov2 [17], ClueWeb09 [1], and CC-News [38]. For Gov2 and ClueWeb09 collections, we used the freely-available inverted indexes and query logs by Daniel Lemire (see [34] for details), corresponding to the URL-sorted document enumeration [48] (which tends to yield runs of successive elements in the sets). The query log contains 20,000 random queries from the TREC million-query track (1MQ). Each query has at least 2 query terms. Also, each term is in the top-1M most frequently queried terms. For CC-News we use the freely-available inverted index by Mackenzie et al. [38] in Common Index File Format (CIFF) [37], as well as their query log of 9,666 queries. Table 1 shows a summary of statistics of the collections. In all cases, we only keep sets with at least 4,096 elements.

| Table 1 Dataset summary and average space usage (in bits per integer, bpi) for different compression measures and baseline representations. |
|---|---|---|
| Gov2 | ClueWeb09 | CC-News |
| # Lists | 57,225 | 131,567 | 79,831 |
| # Integers | 5,509,206,378 | 14,895,136,282 | 18,415,151,585 |
| $u$ | 25,205,179 | 50,220,423 | 43,495,426 |
| $\lceil \lg u \rceil$ | 25 | 26 | 26 |
| gap($S$) | 2.25 | 3.25 | 3.70 |
| rle($S$) | 1.99 | 3.33 | 4.23 |
| trie($S$) | 3.48 | 4.56 | 5.18 |
| rTrie($S$) | 2.51 | 4.00 | 5.12 |
| Elias $\gamma$ | 3.71 | 5.74 | 6.81 |
| Elias $\delta$ | 3.64 | 5.40 | 6.69 |
| Fibonacci | 3.90 | 5.35 | 6.09 |
| Elias $\gamma$ 128 | 4.07 | 6.10 | 7.05 |
| Elias $\delta$ 128 | 4.00 | 5.77 | 7.17 |
| Fibonacci 128 | 4.26 | 5.71 | 6.45 |
| rrr_vector<> | 11.82 | 19.94 | 11.29 |
| sd_vector<> | 8.45 | 8.52 | 7.17 |
As baseline, Table 1 also shows the average bit per integer (bpi) for different compression measures on our tested set collections. We also show the average bpi for different integer compression approaches, namely Elias $\gamma$ and $\delta$ [22], Fibonacci [25], $r_{rrr\text{\_vector}}$ [47], and $s_{sd\text{\_vector}}$ [42], all of them from the sdsl library [27]. In particular, Elias $\gamma$, $\delta$, and Fibonacci codes are known for yielding highly space-efficient set representations in IR indexing [15], hence they are a strong baselines for comparison. We show a plain version of them, as well as variants with blocks of 128 integers. The latter are needed to speed up decoding. However, these approaches are relatively slow to be decoded [15, See Table 6.9], and hence yield higher intersection times. On the other hand, $s_{sd\text{\_vector}}$ uses $n \log (u/n) + 2n + o(n)$ bits to encode a set of $n$ elements and universe $[0..u]$. Finally, $r_{rrr\text{\_vector}}$ uses $B(n, u) + o(u)$ bits of space. As it can be seen, the $o(u)$-bit term yields a higher space usage.

Next, we compare our approaches with state-of-the-art set compression alternatives available at the project Performant Indexes and Search for Academia 1 (PISA) [39]:
- IPC: the Binary Interpolative Coding approach by Moffat et al. [40]. This is a highly space-efficient approach, with a relatively slow processing performance [15, 40].
- PEF Opt: the highly competitive approach by Ottaviano and Venturini [43].
- OptPFD: The Optimized PForDelta approach by Yan et al. [53].
- SIMD-BP128: The highly efficient approach by Lemire and Boytsov [35], aimed at decoding billions of integers per second using vectorization capabilities of modern processors.
- Simple16: The approach by Zhang et al. [54], a variant of the Simple9 approach [5] that combines a relatively good space usage and an efficient intersection time.
- VarintGB: The approach used in Google and presented by Dean [19].
- Varint-G8IU: by Stepanov et al. [49], using SIMD instructions to speed-up set processing.

We also compared with the following approaches, available from their authors:
- Roaring: the compressed bitmap approach by Lemire et al. [36], widely used as indexing tool on several systems and platforms [3]. Roaring bitmaps are highly competitive, leveraging modern CPU hardware architectures. We use the code from the authors [2].
- RUP: The recent recursive universe partitioning approach by Pibiri [45], using also SIMD instructions to speed up processing. We use the code from the author [44].

Table 2 shows the average experimental intersection time (in milliseconds per query) and space usage (in bits per integer) for all the alternatives tested. Figure 7 (in the Appendix) shows the same results, using space vs. time plots. Our approaches introduce competitive trade-offs, as follows:

**Results for Gov2:** $r_{Trie}$ uses 1.16–1.329 times the space of PEF, the former being 1.549–2.442 times faster. $r_{Trie}$ uses 0.481–0.548 times the space of Roaring, the former being up to 1.415 times faster. Finally, $r_{Trie}$ uses 0.837–0.954 times the space of RUP, the former being up to 1.428 times faster.

**Results for ClueWeb09:** $r_{Trie}$ uses 1.188–1.361 times the space of PEF, the former being 2.117–3.316 times faster. Also, $r_{Trie}$ uses 0.551–0.631 times the space of Roaring, the former being 1.221–1.913 times faster. Finally, $r_{Trie}$ uses 0.823–0.943 times the space of RUP, the former being 1.391–2.178 times faster.

**Results for CC-News:** for this dataset, the resulting inverted lists have considerably less runs, hence the space usage of trie and $r_{Trie}$ are about the same. However, trie is faster than $r_{Trie}$, as the code to handle runs introduces an overhead that does not pay

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1 https://github.com/pisa-engine/pisa
off in this case. So, we will use \textit{trie} to compare here. It uses 1.512–1.722 times the space of \textit{PEF}, the former being 1.987–3.326 times faster. \textit{trie} uses 0.889–1.013 times the space of \textit{Roaring}, the former being up to 1.067 times faster. Finally, \textit{trie} uses 1.043–1.188 times the space of \textit{RUP}, the former being up to 1.044 times faster.

We can conclude that in all tested datasets, at least one of our trade-offs is the fastest and competitive in space usage, outperforming the highly-engineered ultra-efficient set compression techniques we tested.

\section{Conclusions}

Trie partition certificates, the main concept we introduced as an alternative to existing certificates by Demaine et al.\cite{20} and Barbay and Kenyon\cite{12}, allowed us to introduce our main contributions. In particular, we were able to prove that Trabb-Pardo’s intersection algorithm\cite{50} works in $O(k \delta \log(u/\delta))$ time, where $\delta$ is the alternation measure of the query instance\cite{12}. Thus, Trabb-Pardo’s intersection algorithm was likely the first adaptive intersection algorithm that ever existed, appearing about 22 years before Demaine et al.’s adaptive approach. The lack of analysis on this algorithm (the original author only analyzed his algorithm in the average case) might explain the lack of consideration regarding this algorithm, in particular in practice. Motivated by this result, we introduced compressed representations of integer sets preserving the running time of Trabb-Pardo’s algorithm, and even improving it. Summarizing, our proposals: (1) use compressed space usage, (2) have adaptive intersection computation time, and (3) have highly competitive practical performance.

Multiple avenues for future research are open now. For instance, novel data structures supporting operation $\text{rank}_1$ have emerged recently\cite{32}. These offer interesting trade-offs, using less space than then ones we used, with competitive operation times. Another interesting
line is that of alternative binary trie compact representations. E.g., a DFS representation [13]
(rather than BFS, as the one used in this paper), which would potentially reduce the number
of cache misses when traversing the tries. Finally, our representation would support dynamic
sets (where insertion and deletion of elements are allowed) if we use dynamic binary tries [6].

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Trie-Compressed Adaptive Set Intersection


A Plots of Experimental Results

**Figure 7** Space vs. time trade-off for all alternative tested on the 3 datasets.
Approximation Algorithms for the Longest Run Subsequence Problem

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Abstract

We study the approximability of the Longest Run Subsequence problem (LRS for short). For a string \(S = s_1 \cdots s_n\) over an alphabet \(\Sigma\), a run of a symbol \(\sigma \in \Sigma\) in \(S\) is a maximal substring of consecutive occurrences of \(\sigma\). A run subsequence \(S'\) of \(S\) is a sequence in which every symbol \(\sigma \in \Sigma\) occurs in at most one run. Given a string \(S\), the goal of LRS is to find a longest run subsequence \(S^*\) of \(S\) such that the length \(|S^*|\) is maximized over all the run subsequences of \(S\). It is known that LRS is APX-hard even if each symbol has at most two occurrences in the input string, and that LRS admits a polynomial-time \(k\)-approximation algorithm if the number of occurrences of every symbol in the input string is bounded by \(k\). In this paper, we design a polynomial-time \(k+1\) approximation algorithm for LRS under the \(k\)-occurrence constraint on input strings. For the case \(k = 2\), we further improve the approximation ratio from \(\frac{3}{2}\) to \(\frac{4}{3}\).

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1 Introduction

The main goal of genome analysis is to study and compare genetic content among organisms, and thus genome sequencing to determine the complete sequence of a genome is one of its most important stages. Since the first whole genome was obtained [10], genome sequencing technologies have significantly improved. Almost all the current DNA sequencing technologies are based on the following process: First, tens or hundreds of millions of fragments from random positions on the DNA sequence are read via shotgun sequencing. Second, these randomly extracted fragments, called reads, are merged to form a set of contiguous sequences, called contigs, by using an assembly algorithm. Then, the contigs are ordered correctly in a phase called scaffolding. One commonly used approach for scaffolding is to rearrange contigs by comparing two or more incomplete assemblies of related samples (see, for example, [8]).
Approximation Algorithms for the Longest Run Subsequence Problem

In the context of the scaffolding phase of genome assembly, the One-Sided Scaffold Filling problem [9], Two-Sided Scaffold Filling problem [7], One-Side-Filled Longest Common Subsequence problem [3], and Two-Side-Filled Longest Common Subsequence problem [4] were formulated as combinatorial optimization problems on two strings. For those problems, their computational complexities were proved, and then fixed-parameter tractable algorithms, approximation algorithms, and exponential-time exact algorithms were proposed in [2, 3, 4, 7]. Very recently, as a different formulation of the scaffolding phase, Schrinner et al. [11, 12] introduced the Longest Run Subsequence problem (LRS for short), defined as follows: For a string $S = s_1 \cdots s_n$ over an alphabet $\Sigma$, a run of a symbol $\sigma \in \Sigma$ in $S$ is a maximal substring of consecutive occurrences of $\sigma$. A run subsequence $S'$ of $S$ is a sequence in which every symbol $\sigma \in \Sigma$ occurs in at most one run. Given a string $S$, the goal of LRS is to find a longest run subsequence $S^*$ of $S$ such that the length $|S^*|$ is maximized over all the run subsequences of $S$.

**Example 1.** Consider the string $S = abacacbab$ over the alphabet $\Sigma = \{a, b, c\}$. It contains (i) four runs of symbol $a$, i.e., $a$ in the first position, $a$ in the third position, $a$ in the fifth position, and $a$ in the ninth position, (ii) three runs of symbol $b$, i.e., $b$ in the second position, $bb$ in the seventh and eighth positions, and $b$ in the tenth position, and (iii) two runs of $c$, i.e., $c$ in the fourth position, and $c$ in the sixth position in $S$. The numbers of occurrences of $a$, $b$, and $c$ are four, four, and two, respectively.

An optimal solution to LRS on input $S$ is $S^* = aaccbbb$. For example, the leftmost run $aa$ of length two in $S^*$ is obtained from the leftmost substring $a\text{bc}a$ in $S$ by deleting the second character $b$. One sees that $S^*$ is a run subsequence, i.e., $S^*$ contains (at most) one run for every symbol. The length of $S^*$ is seven. Note that $S' = a\text{acacb}$ is another optimal solution since $|S'|$ is also seven.

Schrinner et al. [12] showed that LRS is NP-hard. Subsequently, Dondi and Sikora [5] showed that LRS is APX-hard even if each symbol has at most two occurrences in the input string, and that LRS admits a polynomial-time $\min\{|\Sigma|, k\}$-approximation algorithm if the number of occurrences of every symbol in the input string is bounded by $k$.

In this paper, we propose the following improved approximation algorithms for LRS:

- For the case $k = 2$, we further improve the approximation ratio from $\frac{3}{2}$ to $\frac{4}{3}$.

**Related work.** The fixed-parameter tractability and the parameterized complexity of LRS have been previously investigated [5, 12]: Schrinner et al. [12] showed that there is an $O(|\Sigma| \cdot |S| \cdot 2^{|\Sigma|})$-time algorithm, given a string $S$ over an alphabet $\Sigma$ as input of LRS, i.e., LRS is fixed-parameter tractable when parameterized by the size $|\Sigma|$ of the alphabet on which the input string is defined. Dondi and Sikora [5] showed that LRS can be solved by a randomized algorithm in $O(2^r \cdot r \cdot |S|^3)$ time and polynomial space, where $r$ is the number of different runs in a solution, and thus $r \leq |S|$. They also proved that LRS admits a polynomial kernel when parameterized by the length of the solution, but that it does not admit a polynomial kernel when parameterized by the size $|\Sigma|$ of the alphabet or by the number $r$ of runs.

2 Preliminaries

Let $\Sigma$ be a finite alphabet of symbols. A string $S = s_1 \cdots s_n$ is a sequence of $n$ characters, each of which is a symbol in $\Sigma$. Two or more characters in $S$ can be the same symbol in $\Sigma$. For a string $S = s_1 \cdots s_n$, $|S|$ denotes the length of $S$, i.e., $|S| = n$. A subsequence of $S$ is a
sequence $s_1, \cdots, s_m$, such that $1 \leq i_1 < i_2 < \cdots < i_m \leq |S|$. Let $S[i]$ denote the character of $S$ in the $i$th position for $1 \leq i \leq |S|$, and $S[i, j]$ denote the substring of $S$ that starts from the $i$th position and ends at the $j$th position. For a symbol $\sigma$, we denote by $\sigma^k$ a string that is the concatenation of $k$ occurrences of symbol $\sigma$ for some integer $k \geq 1$. A run in $S$ is a substring $S[i, j]$ such that: (1) $S[i] = S[i + 1] = \cdots = S[j]$; (2) $S[i - 1] \neq S[i]$ if $i > 1$; and (3) $S[j + 1] \neq S[j]$ if $j < |S|$. For any $\sigma \in \Sigma$, a run in $S$ of the form $\sigma^k$ is called a length-$k$ $\sigma$-run in $S$. Observe that if $S[i, j]$ is a $\sigma$-run, then it has length $j - i + 1$. Given a string $S$ on alphabet $\Sigma$, a run subsequence $S'$ of $S$ is a subsequence in which every symbol $\sigma \in \Sigma$ occurs in at most one run.

Let $\text{occ}(\sigma)$ be the number of occurrences of $\sigma$ in the input string $S$. Let $\text{occ}_{\text{max}}(S) = \max_{\sigma \in S} \text{occ}(\sigma)$. For example, consider a string $S = abaceabab$. Then, $S$ includes four $a$-runs, $a$, $a$, $a^2$, and $a$, three $b$-runs, $b$, $b^2$, and $b$, and one length-1 $c$-run. The number $\text{occ}(a)$ of occurrences of $a$ is five. Also, $\text{occ}(b) = 4$ and $\text{occ}(c) = 1$. Therefore, $\text{occ}_{\text{max}}(S) = 5$.

Our problem LRS can be formulated as follows:

\begin{itemize}
    \item \textbf{Problem 2 (Longest Run Subsequence problem, LRS).} Given an alphabet $\Sigma$ and a string $S = s_1 \cdots s_n$ with $s_i \in \Sigma$, the goal of LRS is to find a longest run subsequence $S^*$ of $S$, i.e., every $\sigma \in \Sigma$ occurs in at most one run in $S^*$ and the length $|S^*|$ is maximized over all the run subsequences of $S$.
\end{itemize}

Schirrner et al. [12] show that LRS is NP-hard by giving a polynomial-time reduction from the \textsc{Linear Ordering Problem}, which is shown to be NP-hard in [6]. In this paper we consider the following restricted LRS:

\begin{itemize}
    \item \textbf{Problem 3 ($k$-Longest Run Subsequence problem, $k$-LRS).} If the maximum number $\text{occ}_{\text{max}}(S)$ of occurrences of symbols in the input $S$ is bounded by $k$, then the problem is called the $k$-Longest Run Subsequence problem, $k$-LRS.
\end{itemize}

One sees that 1-LRS is trivial since the length of all the runs in the input string $S$ is one, and thus the input $S$ itself is the optimal run subsequence. Dondi and Sikora [5] show that 2-LRS remains hard even from the approximation point of view; they give an L-reduction from the \textsc{Minimum Independent Set on Cubic Graph Problem}, which is shown to be APX-hard in [1]:

\begin{itemize}
    \item \textbf{Proposition 4 ([5]).} 2-LRS is APX-hard.
\end{itemize}

Suppose that an input string of $k$-LRS is $S$ over an alphabet $\Sigma$. Also, without loss of generality, we assume here that every symbol in $\Sigma$ appears at least once, and the maximum number $\text{occ}_{\text{max}}(S)$ of occurrences of symbols in $S$ is $k$. Note that the length of an optimal run subsequence is bounded by $k|\Sigma|$. Consider the following two simple algorithms, (i) and (ii):

\begin{itemize}
    \item (i) Arbitrarily select one run of every symbol $\sigma \in \Sigma$ in $S$, and construct a run subsequence $S'$ by concatenating all the selected runs.
    
    One sees that $|S'|$ is at least $|\Sigma|$. Therefore, we can conclude that $k$-LRS is $k$-approximable.

    \item (ii) Find a symbol, say, $\sigma$ of the maximum occurrences $k$, and construct another run subsequence $S'' = \sigma^k$.
    
    Then, we can conclude that $k$-LRS is $|\Sigma|$-approximable. By using those two algorithms, we obtain the following proposition:
\end{itemize}

\begin{itemize}
    \item \textbf{Proposition 5 ([5]).} There is a min$(|\Sigma|, k)$-approximation algorithm for $k$-LRS.
\end{itemize}

Since min$(|\Sigma|, k) \leq \sqrt{|S|}$, the above proposition implies the following:

\begin{itemize}
    \item \textbf{Corollary 6 ([5]).} The general LRS problem admits a $\sqrt{|\Sigma|}$-approximation algorithm.
\end{itemize}
In this section, we improve the approximation ratio for $k$-LRS from $k$ to $\frac{k+1}{2}$. Our approximation algorithm $\text{ALG}$ uses a very natural idea:

Algorithm $\text{ALG}$. Given an input string $S$ over an alphabet $\Sigma$, $\text{ALG}$ selects a longest $\sigma$-run in $S$ for each $\sigma \in \Sigma$, and outputs the concatenation of all the selected longest runs.

Example 7. Consider the input string $S = abacaabbab$ (for 5-LRS). The longest $a$-run, $b$-run, $c$-run are $aa$ in the fifth and sixth positions, $bb$ in the seventh and eighth positions, and $c$ in the fourth position. Therefore, the output of $\text{ALG}$ is $\text{ALG} = caabb$. \hfill \text{∎}

We now prove that the above simple algorithm achieves the claimed approximability bound:

Theorem 8. $\text{ALG}$ is a polynomial-time $\frac{k+1}{2}$-approximation algorithm for $k$-LRS.

Proof. Clearly, $\text{ALG}$ returns a valid solution since one run is selected for every symbol in $S$, and runs in polynomial time. We bound its approximation ratio in the following. Let $S$ be an input string of $k$-LRS. We assume that $S$ consists of $m$ symbols, i.e., $|\Sigma| = m$, and $\text{occ}_{\text{max}}(S) = k$. Then, suppose that $\text{OPT}$ and $\text{ALG}$ are solutions obtained by an optimal algorithm and our algorithm $\text{ALG}$, respectively, for the input $S$. We consider the following two cases: (Case 1) The length of every run in $S$ is one, and (Case 2) the length of some run in $S$ is at least two.

(Case 1). Suppose that the length of every run in $S$ is one. Let $m_\ell$ be the number of symbols in $\text{OPT}$ such that the length of the run of those symbols is exactly $\ell$ ($\leq k$).

First, the following two equalities hold:

$$|\text{OPT}| = \sum_{i=1}^{k} i \cdot m_i; \text{ and}$$

$$|\text{ALG}| = m. \quad (1)$$

Let $D$ be the number of characters deleted from $S$ by the optimal algorithm. Since $|\Sigma| = \sum_{i=0}^{k} m_i = m$ and $\text{occ}_{\text{max}}(S) = k$, the following is satisfied:

$$|\text{OPT}| = |S| - D \leq km - D. \quad (2)$$

We now derive a lower bound on $D$. Suppose that a symbol $\sigma_2$ in $S$ appears exactly twice in the optimal solution $\text{OPT}$, i.e., $\text{OPT}$ contains the length-2 $\sigma_2$-run $\sigma_2\sigma_2$. Recall that the length of all the runs in the input string $S$ is one. Namely, there is at least one different character, say, $\sigma'$ between two $\sigma_2$’s in $S$. That is, $\sigma'$ must be deleted from $S$ in order to obtain the length-2 $\sigma_2$-run. Since $\text{OPT}$ contains $m_2$ symbols such that the length of the runs of those symbols is exactly two, the total number of deleted characters from $S$ to obtain the length-2 runs is at least $m_2$. It is important to note that the character-deletion to obtain each run is independently carried out, and therefore the number of deleted characters is not doubly counted. Similarly, the total number of deleted characters from $S$ to obtain the length-2 runs is at least $(\ell - 1)m_\ell$ for each $3 \leq \ell \leq k$. As a result, we obtain the following lower bound on $D$:

$$D \geq m_2 + 2m_3 + \cdots + (k-1)m_k = \sum_{i=2}^{k} (i-1)m_i = \sum_{i=1}^{k} (i-1)m_i. \quad (4)$$
From Eq.(3) and Eq.(4), the following inequality holds:

\[ |OPT| \leq km - \sum_{i=1}^{k} (i-1)m_i. \]

From Eq.(1), this can be rewritten as:

\[ |OPT| \leq (k+1)m - |OPT|, \]
and then rearranged to give:

\[ |OPT| \leq \frac{(k+1)m}{2}. \]

From Eq.(2), we obtain the following approximation ratio:

\[ \frac{|OPT|}{|ALG|} \leq \frac{k+1}{2}. \]

(Case 2). Suppose that the length of a \( \sigma \)-run in \( S \) is at least two and \( S \) consists of symbols in \( \Sigma \). For every such symbol \( \sigma \in \Sigma \), we consider a different symbol \( \bar{\sigma} \), called a dummy symbol. Then, we insert \( \bar{\sigma} \) between every consecutive two symbols \( \sigma\sigma \) in \( S \) so that the two \( \sigma \)'s are not consecutive. Hence we obtain a longer sequence \( S_d \) such that the length of all the runs in \( S_d \) is one. For example, consider a string

\[ S = abacaabbab. \]

Then, we insert a dummy \( \bar{\sigma} \) between the fifth and the sixth positions, a dummy \( \bar{b} \) between the seventh and the eighth positions, and the other dummy \( \bar{b} \) between the eighth and the ninth positions as follows:

\[ S_d = abaca\bar{\sigma}ab\bar{b}ab. \]

Note that the number of occurrences of each dummy \( \bar{\sigma} \) is at most \( k-1 \) since the maximum number \( occ_{\max}(S) \) of occurrences of (original) symbols in \( S \) is bounded by \( k \). Suppose that \( OPT_d \) and \( ALG_d \) are solutions obtained by an optimal algorithm and our algorithm \( ALG \), respectively, for the input \( S_d \). One sees that the maximum number \( occ_{\max}(S_d) \) of occurrences of symbols in \( S_d \) is also bounded by \( k \). Therefore, from the arguments in (Case 1), the following inequality is satisfied:

\[ \frac{|OPT_d|}{|ALG_d|} \leq \frac{k+1}{2}. \] (5)

The original input \( S \) is a subsequence of \( S_d \). Hence, the following clearly holds:

\[ |OPT| \leq |OPT_d|. \] (6)

Now consider \( ALG \) and \( ALG_d \). (i) For each symbol \( \sigma \) such that the length of all the \( \sigma \)-runs is one, its dummy \( \bar{\sigma} \) is not inserted into \( S_d \). Hence, \( ALG \) and \( ALG_d \) contain one \( \sigma \), but, of course, neither contains any \( \bar{\sigma} \). (ii) If the maximum length of a \( \sigma \)-run in \( S \) is (at least) two for some symbol \( \sigma \), then \( ALG \) contains (at least) two \( \sigma \)'s. On the other hand, \( ALG_d \) contains one \( \sigma \) and one dummy \( \bar{\sigma} \) instead. From (i) and (ii), we have:

\[ |ALG| \geq |ALG_d|. \] (7)
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From the three inequalities (5), (6), and (7), the following approximation ratio is obtained again:

\[
|OPT| \leq \frac{|OPT_d|}{|ALG_d|} \leq \frac{k+1}{2}.
\]

For both cases (Case 1) and (Case 2), the approximation ratio of \( ALG \) is bounded above by \( \frac{k+1}{2} \).

\[\preceq\]

\[\triangleright\]

Remark 9. To see that the approximation analysis above is tight, consider the following string \( S \), where \( |S| = n = 2k\ell \), and \( \sigma_i \neq \sigma_j \) for \( i \neq j \).

\[
S = \sigma_1\sigma_2\sigma_1\sigma_2 \cdots \sigma_1\sigma_2\sigma_3\sigma_4\sigma_1\sigma_2 \cdots \sigma_{2\ell-1}\sigma_{2\ell-1}\sigma_{2\ell} \cdots \sigma_{2\ell-1}\sigma_{2\ell}.
\]

Namely, the length-2k prefix string contains \( k \sigma_1 \)'s and \( k \sigma_2 \)'s alternatively. The next string of length 2k contains \( k \sigma_3 \)'s and \( k \sigma_4 \)'s alternatively, and so on. Then, we can obtain the following run subsequence \( S' \) by deleting \( k-1 \sigma_2 \)'s from the first length-2k prefix string, \( k-1 \sigma_4 \)'s from the next string of length 2k, and so on:

\[
S' = \sigma_1^k\sigma_2^k\sigma_4^k \cdots \sigma_{2\ell-1}^k\sigma_{2\ell}.
\]

Hence, the length of OPT is at least \( |S'| = (k+1)\ell \). On the other hand, the solution ALG of our algorithm ALG for \( S \) contains one of the \( k \sigma_i \)'s for each \( 1 \leq i \leq 2\ell \):

\[
ALG = \sigma_1\sigma_2 \cdots \sigma_{2\ell}.
\]

The length of ALG is 2\ell. As a result,

\[
\frac{|OPT|}{|ALG|} \geq \frac{k+1}{2}.
\]

This shows that the analysis of the approximation ratio in the proof of Theorem 8 is tight. \[\triangleright\]

Recall that we can always return a run subsequence of length \( k \) as shown in the previous section, and k-LRS is \( |\Sigma| \)-approximable. Therefore, we obtain the following corollary:

\[\triangleright\]

Corollary 10. There is a polynomial-time \( \min\{(|\Sigma|, \frac{k+1}{2}) \}-approximation algorithm for k-LRS.

4 A polynomial-time \( \frac{4}{3} \)-approximation algorithm for 2-LRS

For 2-LRS, ALG achieves the approximation ratio of \( \frac{3}{2} \). In this section we improve the approximation ratio to \( \frac{4}{3} \).

As shown in Remark 9, the following string \( S \) is a bad example for ALG.

\[
S = ababcdcdefef.
\]

One sees that from the leftmost substring \( S[1,4] = abab \) of length four (resp. \( S[5,8] = cced \) and \( S[9,12] = efef \)), we can only obtain a run subsequence of length at most three, i.e., the length of any optimal solution is at most nine. Therefore, one of the possible optimal solution OPT for \( S \) is:

\[
OPT = aabccccdef.
\]
The solution $ALG$ of $ALG$ for $S$ is:

$$ALG = abcdef.$$ 

Namely, $OPT$ has two $a$’s (resp. two $c$’s and two $e$’s), but $ALG$ has only one $a$ (resp. one $c$ and one $e$). This observation suggests to us that if there is only one character, say, $\sigma'$ between two occurrences of a symbol $\sigma$, then we should delete $\sigma'$ and obtain a run $\sigma\sigma$ of length two. This is a basic strategy of our new algorithm $ALG_2$.

Before describing details of $ALG_2$, we give some definitions which are used in the following. Let $S$ be an input string. Assume that all the symbols in $\Sigma$ appear in $S$. We define several subsets of $\Sigma$ in the following.

- Let $\Sigma_1 = \{ \sigma \mid \text{occ}(\sigma) = 1, \sigma \in \Sigma \}$ be a set of symbols that appear exactly once in the input string $S$.
- Let $\Sigma_2 = \{ \sigma \mid \text{occ}(\sigma) = 2, \sigma \in \Sigma \}$ be a set of symbols that appear exactly twice in the input string $S$.

Note that $\Sigma = \Sigma_1 \cup \Sigma_2$ in 2-LRS. Now, we consider a symbol $\sigma \in \Sigma_2$ and define several disjoint subsets of $\Sigma_2$. In the following, by distance we mean the number of characters between the two occurrences of a symbol.

- If two $\sigma$’s consecutively appear in $S$, then we call $\sigma$ a distance-0 symbol. Let $\Sigma_{2,0}$ be a subset of all the distance-0 symbols in $\Sigma_2$.
- If there is one character between two $\sigma$’s, then we call $\sigma$ a distance-1 symbol. Let $\Sigma_{2,1}$ be a subset of all the distance-1 symbols in $\Sigma_2$.
- We define $\Sigma_{2,\geq 2} = \Sigma_2 \setminus (\Sigma_{2,0} \cup \Sigma_{2,1})$, i.e., for each $\sigma \in \Sigma_{2,\geq 2}$, $\sigma$ appears twice in $S$ and there are at least two characters between the two $\sigma$’s.

Next, consider a symbol $\gamma \in \Sigma_1$. As a special case, the left and the right symbols of $\gamma$ can be the same symbol $\gamma’ \in \Sigma_{2,1}$, i.e., the input string $S$ possibly contains a substring $\gamma’\gamma’$ of length 3, called a special triple.

- Let $\Gamma_1$ be a set of center symbols of special triples. Note that $\Gamma_1 \subseteq \Sigma_1$.
- Let $\Gamma_{2,1}$ be a set of left and right symbols of special triples. Note that $\Gamma_{2,1} \subseteq \Sigma_{2,1}$.

One sees that $|\Gamma_1| = |\Gamma_{2,1}|$.

Finally, consider two symbols $\sigma$ and $\sigma'$ in $\Sigma_{2,1} \setminus \Gamma_{2,1}$ in the input string $S$ such that the substring(s) containing $\sigma$ and $\sigma'$ can be represented by (i) $S = \cdots \sigma'\sigma' \cdots$, or (ii) $S = \cdots \sigma \lambda \sigma \cdots \sigma' \lambda' \sigma' \cdots$, where both $\lambda$ and $\lambda'$ are in $\Sigma_{2,\geq 2}$. (i) If $S$ contains $\sigma'\sigma'\sigma'$ as a substring, then we say that a pair of $\sigma$ and $\sigma'$ is called a $\Psi$-pair. Then, $\sigma$ and $\sigma'$ belong to a set $\Psi_{2,1}$. (ii) If $\lambda = \lambda'$, then we say that a pair of $\sigma$ and $\sigma'$ is a $\Lambda$-pair related to $\lambda$. Then, $\sigma$ and $\sigma'$ belong to a set $\Lambda_{2,1}$ and $\lambda$ belongs to $\Lambda_{2,\geq 2}$. Note that $|\Lambda_{2,1}| = 2|\Lambda_{2,\geq 2}|$.

**Algorithm.** The following is a description of our algorithm $ALG_2$. During execution of $ALG_2$, we determine which characters are included into the run subsequence $ALG_2$ or not, step by step. Finally, $ALG_2$ outputs the concatenation of the characters (or the subsequences) included into $ALG_2$ in each step.

**Algorithm $ALG_2$.**

- **Input** An input string $S$ over an alphabet $\Sigma$ such that every symbol in $\Sigma$ appears at most twice.
- **Output** A run subsequence.
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1. Count the number of occurrences of every symbol in \( \Sigma \), and divide \( \Sigma \) to two subsets \( \Sigma_1 \) and \( \Sigma_2 \). Then, examine the distance of every symbol in \( \Sigma_2 \), and obtain \( \Sigma_{2,0}, \Sigma_{2,1}, \text{ and } \Sigma_{2,2} \).

2. Find all the special triples, all the \( \Psi \)-pairs, and all the \( \Lambda \)-pairs.

3. For every \( \sigma \in \Sigma_{2,0} \), the length-2 \( \sigma \)-run \( \sigma^2 \) is included into \( ALG_2 \).

4. For every \( \sigma \in \Sigma_{2,1} \), execute the following:
   a. For every special triple \( \gamma' \gamma' \), the first two characters \( \gamma \in \Gamma_{2,1} \) and \( \gamma \in \Gamma_1 \) are included into \( ALG_2 \). That is, the third character \( \gamma' \) of that special triple is not included into \( ALG_2 \).
   b. For every \( \Psi \)-pair of \( \sigma \) and \( \sigma' \), i.e., for each string \( \sigma \sigma' \sigma' \), its subsequence \( \sigma \sigma' \sigma' \) is included into \( ALG_2 \). That is, the third character \( \sigma \) of that string is not included into \( ALG_2 \).
   c. For every \( \Lambda \)-pair related to \( \lambda \) of \( \sigma \) and \( \sigma' \), i.e., for two strings \( \sigma \lambda \sigma \) and \( \sigma' \lambda \sigma' \), two subsequences \( \sigma \lambda \), and \( \sigma' \lambda \sigma' \) are included into \( ALG_2 \). That is, the third character \( \lambda \) of the latter string are not included into \( ALG_2 \).
   d. For every \( \sigma \in \Sigma_{2,1} \setminus (\Gamma_{2,1} \cup \Psi_{2,1} \cup \Lambda_{2,1}) \), \( \sigma^2 \) is included into \( ALG_2 \). That is, the character between the two \( \sigma \)'s is not included into \( ALG_2 \).

5. For every \( \sigma \in \Sigma_{2,2} \setminus \Lambda_{2,2} \), only the first occurrence of \( \sigma \) is included into \( ALG_2 \). That is, if neither of the two occurrences of \( \sigma \) is determined whether or not to be included into \( ALG_2 \), then the first occurrence is included into \( ALG_2 \) and the other not into \( ALG_2 \). If only one occurrence remains undetermined, then it is included into \( ALG_2 \).

6. Every \( \sigma \in \Sigma_1 \setminus \Gamma_1 \) is included into \( ALG_2 \).

7. Output the concatenation of the characters and the subsequences included into \( ALG_2 \) in Step 3 through Step 6 as a run subsequence, and then halt.

\[ \text{Step 4.} \]
\[ \text{Step 5.} \]
\[ \text{Step 6.} \]

---

1. Importantly, the output run subsequence of \( ALG_2 \) includes at least one occurrence of every symbol in \( \Sigma \).

---

2. To clarify the behavior of \( ALG_2 \), we take a look at the following input string of length 20:

\[ S = abacdbdecefgfhhijjk. \]

One sees that \( \Sigma_1 = \{g, i\}, \Sigma_{2,0} = \{h\}, \Sigma_{2,1} = \{a, d, e, f, j, k\}, \text{ and } \Sigma_{2,2} = \{b, c\}. \]

\( S[14, 15] = hh \) is included into \( ALG_2 \). (Step 4-(i)) Since \( f \in \Sigma_{2,1} \) and \( g \in \Sigma_1 \), \( S[10, 12] = fgf \) is a special triple. Therefore, we select \( fg \) from \( fgf \). (Step 4-(ii)) Since there is a substring \( S[17, 20] = jkjk \), the pair of \( j \) and \( k \) is a \( \Psi \)-pair, \( \Psi_{2,1} = \{j, k\} \). Then, \( jkk \) is included into \( ALG_2 \). (Step 4-(iii)) \( S \) contains \( S[1, 3] = aba \) and \( S[5, 7] = dbd \) and thus the pair of \( a \) and \( d \) is a \( \Lambda \)-pair related to \( b \); \( \Lambda_{2,1} = \{a, d\} \) and \( \Lambda_{2,2} = \{b\} \). Hence, \( ab \) and \( dd \) are included into \( ALG_2 \). (Step 4-(iv)) From \( S[8, 10] = cce \), we obtain a run \( c^2 \) of length two, and \( S[9] = c \) is not included into \( ALG_2 \). (Step 5) The fourth character \( c \) is included into \( ALG_2 \) since \( c \in \Sigma_{2,2} \setminus \Gamma_{2,2} \) and \( S[9] = c \) is not included into \( ALG_2 \) in Step 4-(iv). (Step 6) The

---

Alternative, we can choose any one of the two occurrences of each symbol, to obtain the same approximation ratio.
16th character \( i \) is included into \( \text{ALG}_2 \) since \( i \in \Sigma_1 \setminus \Gamma_1 \). (Step 7) Finally, the following concatenation of the characters and the subsequences obtained in Step 3 through Step 6 is output as the run subsequence \( \text{ALG}_2 \) of length 15:

\[
\text{ALG}_2 = abcdde.fghhijklk.
\]

**Theorem 13.** \( \text{ALG}_2 \) is a polynomial-time \( \frac{3}{4} \)-approximation algorithm for 2-LRS.

**Proof.** Clearly, \( \text{ALG}_2 \) returns a valid solution and runs in polynomial time. We bound its approximation ratio in the following. Suppose that \( \text{OPT} \) and \( \text{ALG}_2 \) are run subsequences obtained by an optimal algorithm and our algorithm \( \text{ALG}_2 \), respectively, for the input string \( S \).

We assume that the optimal run subsequence \( \text{OPT} \) consists of the following symbols (OPT1 through OPT4) or characters in special triples (OPT5):

- (OPT1) Consider symbols in \( \Sigma_{2, \geq 2} \). Suppose that there are \( m_{2, \geq 2} \) symbols such that two occurrences of each of them are included into \( \text{OPT} \) by deleting all the characters between two occurrences. Also, suppose that there are \( m_{2, \geq 2} \) symbols such that one occurrence (resp. no occurrence) of each of them is included into \( \text{OPT} \).

- (OPT2) Consider symbols in \( \Sigma_{2, 1} \setminus \Gamma_{2, 1} \). Assume that there are \( m_{2, 1, 2} \) symbols such that two occurrences of each of them are included into \( \text{OPT} \) by deleting one character between two occurrences. Also, suppose that there are \( m_{2, 1, 1} \) symbols such that one occurrence (resp. no occurrence) of each of them is included into \( \text{OPT} \).

- (OPT3) Consider symbols in \( \Sigma_{2, 0} \). Suppose that there are \( m_{2, 0, 2} \) symbols such that two occurrences (resp. no occurrence) of each of them are included into \( \text{OPT} \). Remark that since the goal is to maximize the length of the run subsequence, we can assume that two occurrences (one run of length two) of the symbol in \( \Sigma_{2, 0} \) are completely included into \( \text{OPT} \), or completely deleted.

- (OPT4) Consider symbols in \( \Sigma_1 \setminus \Gamma_1 \). Suppose that there are \( m_{1, 1} \) symbols such that one occurrence (resp. no occurrence) of each of them is included into \( \text{OPT} \).

- (OPT5) Consider special triples. For example, take a look at \( \gamma', \gamma \) where \( \gamma \in \Gamma_1 \) and \( \gamma' \in \Gamma_{2, 1} \). One sees that we cannot select all the three characters into any solution subsequence since it can contain at most one run for every symbol. Therefore, \( \text{OPT} \) includes at most two characters of the special triple, \( \gamma'^2, \gamma', \) or \( \gamma \gamma' \). Since the goal is to maximize the length of the run subsequence, we can assume that \( \text{OPT} \) includes one of the two characters of the special triple, or does not include any character from the special triple. Suppose that there are \( m_{\gamma, 2} \) special triples such that two characters (resp. no character) of each of them are included into \( \text{OPT} \).

Then, the length of \( \text{OPT} \) is calculated as follows:

\[
|\text{OPT}| = 2m_{2, \geq 2} + m_{2, \geq 2} + 2m_{2, 1, 2} + m_{2, 1, 1} + 2m_{2, 0, 2} + m_{1, 1} + 2m_{\gamma, 2}.
\]

Now, let \( D \) be the number of deleted symbols from \( S \) by the optimal algorithm. Then, \( D \) is counted by the above assumption:

\[
D = m_{2, \geq 2} + m_{2, \geq 2} + m_{2, 1, 1} + m_{2, 1, 0} + 2m_{2, 0, 0} + m_{1, 0} + m_{\gamma, 2} + 3m_{\gamma, 0}.
\]
for symbols in $\Sigma_{2,1} \setminus \Gamma_{2,1}$, we assumed in (OPT2) that there are $m_{2,1,2}$ symbols such that two occurrences of each of them are included into $OPT$, i.e., one character between the two occurrences must be deleted. As a result, the following inequality holds:

$$D \geq 2m_{2,2,2} + m_{2,1,2}. \quad (10)$$

Now, we estimate the length of the output run subsequence of $ALG_2$.

(ALG1) Consider symbols in $\Sigma_{2,0}$. In Step 3, two occurrences of every symbol in $\Sigma_{2,0}$ are included into $ALG_2$, i.e., $2m_{2,0,2} + 2m_{2,0,0}$ characters are included into $ALG_2$.

(ALG2) Consider symbols in $\Gamma_{2,1}$. In Step 4-(i), one occurrence of every symbol in $\Gamma_{2,1}$ is included into $ALG_2$, i.e., $m_{\gamma,2} + m_{\gamma,0}$ characters are totally included in $ALG_2$.

(ALG3) Consider symbols in $\Sigma_1$. In Step 4-(i), every symbol in $\Sigma_1$ (i.e., $\Sigma_1 \setminus \Gamma_{2,1}$) is included into $ALG_2$. In Step 6, every symbol in $\Gamma_1$ (i.e., $\Gamma_1 \setminus \Sigma_1$) is included into $ALG_2$. That is, all the symbols in $\Sigma_1$ are included into $ALG_2$. In total, $m_{1,1} + m_{1,0} + m_{\gamma,2} + m_{\gamma,0}$ characters are included into $ALG_2$.

(ALG4) Consider symbols in $\Sigma_{2,2}$. In Step 4-(iii), one occurrence of every symbol in $\Lambda_{2,2}$ (i.e., $\Sigma_{2,2} \setminus \Gamma_{2,1}$) is included into $ALG_2$. Also, in Step 5, one occurrence of every symbol in $\Sigma_{2,2} \setminus \Lambda_{2,2}$ is included into $ALG_2$. In total, $m_{2,2,2} + m_{2,2,1} + m_{2,2,0}$ characters are included into $ALG_2$.

(ALG5) Consider symbols in $\Sigma_{2,1} \setminus \Gamma_{2,1}$. Recall that $|\Sigma_{2,1} \setminus \Gamma_{2,1}| = m_{2,1,2} + m_{2,1,1} + m_{2,1,0}$. Consider a $\Psi$-pair of $\sigma$ and $\sigma'$, i.e., a substring $\sigma \sigma' \sigma'$ of length four in $S$. In Step 4-(ii), three characters $\sigma$, $\sigma'$, and $\sigma''$ are selected from the $\Psi$-pair of $\sigma$ and $\sigma'$. Namely, we can see that three characters per two symbols are included into $ALG_2$. Also, in Step 4-(iii), three characters $\sigma$, $\sigma'$, and $\sigma''$ are selected from every $\Lambda$-pair of $\sigma$ and $\sigma'$. Again, three characters per two symbols are included into $ALG_2$. In Step 4-(iv), two occurrences of every symbol in $(\Sigma_{2,1} \setminus \Gamma_{2,1}) \setminus (\Psi_{2,1} \cup \Lambda_{2,1})$ are included into $ALG_2$. As a result, at least $\frac{3}{2}(m_{2,1,2} + m_{2,1,1} + m_{2,1,0})$ characters are included into $ALG_2$.

Then, the following inequality on the length of $ALG_2$ holds:

$$|ALG_2| \geq m_{2,2,2} + m_{2,2,1} + m_{2,2,0} + \frac{3}{2}(m_{2,1,2} + m_{2,1,1} + m_{2,1,0}) + 2m_{2,0,2} + 2m_{2,0,0} + m_{1,1} + m_{1,0} + 2m_{\gamma,2} + 2m_{\gamma,0}. \quad (11)$$

From Eq.(9) and Eq.(10), we obtain the following inequality:

$$\frac{1}{3}(m_{2,2,1} + 2m_{2,2,0} - m_{2,1,2} + m_{2,1,1} + 2m_{2,1,0} + m_{1,0} + m_{\gamma,2} + 3m_{\gamma,0}) \geq \frac{2}{3}m_{2,2,2}. \quad (12)$$

Therefore, from Eq.(8) and Eq.(12), $|OPT|$ is bounded as follows:

$$|OPT| = \left(\frac{4}{3}m_{2,2,2} + \frac{2}{3}m_{2,1,2} + \frac{2}{3}m_{2,0,2} + m_{2,1,1} + m_{2,0,1} + 2m_{1,1} + 2m_{1,0} + 2m_{\gamma,2} + 2m_{\gamma,0}\right) \leq \left(\frac{4}{3}m_{2,2,2} + \frac{4}{3}m_{2,2,1} + \frac{2}{3}m_{2,2,0} + \frac{5}{3}m_{2,1,2} + \frac{4}{3}m_{2,1,1} + \frac{2}{3}m_{2,1,0} + \frac{2}{3}m_{2,0,2} + \frac{2}{3}m_{2,0,0} + m_{1,1} + \frac{1}{3}m_{1,0} + \frac{7}{3}m_{\gamma,2} + m_{\gamma,0}\right). \quad (13)$$
One can verify that the following is satisfied from Eq.(11) and Eq.(13):
\[
\frac{|OPT|}{|ALG_2|} \leq \frac{4}{3}.
\]

\textbf{Remark 14.} Again, we can show the tightness for the approximation ratio \(\frac{4}{3}\) of \(ALG_2\).

Consider the following string \(S\), where \(|S| = n = 6\ell\).
\[
S = \sigma_1\sigma_2\sigma_3\sigma_4\sigma_5\sigma_6\sigma_4\sigma_5\sigma_6 \cdots \sigma_{3\ell-2}\sigma_{3\ell-1}\sigma_{3\ell-2}\sigma_{3\ell-1}\sigma_{3\ell}.
\]
Then, we can find the following run subsequence \(S'\):
\[
S' = \sigma_1\sigma_2\sigma_3\sigma_4\sigma_5\sigma_6 \cdots \sigma_{3\ell-2}\sigma_{3\ell-1}\sigma_{3\ell}.
\]
Therefore, the length of \(OPT\) is at least \(|S'| = 4\ell\). On the other hand, the solution of our algorithm \(ALG_2\) for \(S\) contains only one of the two \(\sigma_i\)'s for each \(1 \leq i \leq 3\ell\) since every symbol is in \(\Sigma_{2,2}\):
\[
ALG_2 = \sigma_1\sigma_2\cdots\sigma_{3\ell}.
\]
The length of \(ALG_2\) is \(3\ell\). As a result,
\[
\frac{|OPT|}{|ALG_2|} \geq \frac{4}{3}.
\]
This shows that the above approximation analysis is tight.

\section{Conclusion}

We have presented a polynomial-time \(\frac{k+1}{k}\)-approximation algorithm for \(k\)-LRS, where the number of occurrences of every symbol in the input string is at most \(k\). Then, for the case \(k = 2\), we have reduced the approximation ratio to \(\frac{4}{3}\). The current approximation algorithm for 2-LRS is a little bit complicated, and thus might be simplified to obtain the same approximation ratio. Future work is to further improve the approximation ratio of \(\frac{4}{3}\) for 2-LRS, and to design an even better approximation algorithm for general \(k\)-LRS. It would also be useful to derive tight bounds on the polynomial-time approximation hardness of \(k\)-LRS in terms of \(k\).

\section*{References}


Approximation Algorithms for the Longest Run Subsequence Problem


Optimal LZ-End Parsing Is Hard

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Abstract

LZ-End is a variant of the well-known Lempel-Ziv parsing family such that each phrase of the parsing has a previous occurrence, with the additional constraint that the previous occurrence must end at the end of a previous phrase. LZ-End was initially proposed as a greedy parsing, where each phrase is determined greedily from left to right, as the longest factor that satisfies the above constraint [Kreft & Navarro, 2010]. In this work, we consider an optimal LZ-End parsing that has the minimum number of phrases in such parsings. We show that a decision version of computing the optimal LZ-End parsing is NP-complete by showing a reduction from the vertex cover problem. Moreover, we give a MAX-SAT formulation for the optimal LZ-End parsing adapting an approach for computing various NP-hard repetitiveness measures recently presented by [Bannai et al., 2022]. We also consider the approximation ratio of the size of greedy LZ-End parsing to the size of the optimal LZ-End parsing, and give a lower bound of the ratio which asymptotically approaches 2.

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1 Introduction

In the context of lossless data compression, various repetitiveness measures – especially those based on dictionary compression algorithms – and relations between them have recently received much attention (see the excellent survey by Navarro [12, 13]). One of the most fundamental and well-known measures is the LZ77 parsing [15], in which a string is parsed into \( z \) phrases such that each phrase is a single symbol, or the longest substring which has a previous occurrence. LZ-End [9, 10] is a variant of LZ77 parsing with the added constraint that a previous occurrence of the phrase must end at the end of a previous phrase. More formally, the LZ-End parsing is a factorization \( q_1, \ldots, q_z \) of a given string that can be greedily obtained from left to right: each phrase \( q_i \) is either (1) a symbol that is the leftmost occurrence of the symbol or (2) the longest prefix of the remaining suffix \( q_i \cdots q_z \) that is a suffix of \( q_1 \cdots q_j \) for some \( j < i \). It is known that LZ-End parsing can be computed in linear time [6], and there exists a space-efficient algorithm [5].

While there is no known data structure of \( O(z) \) size that provides efficient random access to arbitrary positions in the string, it was recently shown that \( \tilde{O}(1) \) time access could be achieved with \( O(z_e) \) space [8]. Furthermore, concerning the difference between \( z \) and \( z_e \), an upper bound of \( z_e = O(z \log^2(n/z)) \) was shown [8], where \( n \) is the length of the (uncompressed) string. On the other hand, there is an obvious bound of \( z_e = \Omega(z \log n) \) for the unary string, since a previous occurrence of an LZ-End phrase cannot be self-referencing, i.e., overlap with itself, while an LZ77 phrase can. Notice that \( z \leq z_{no} \leq z_e \) holds for any string, where \( z_{no} \) is the number of phrases in the LZ77 parsing that does not allow self-referencing. A family of strings such that the ratio \( z_e/z_{no} \) asymptotically approaches 2 (for large alphabet [10], for binary alphabet [4]) is known, and it is conjectured that \( z_e \leq 2z_{no} \) holds for any strings [10].

While the phrases in the parsings described above are chosen greedily (i.e., longest), we can consider variants which do not impose such constraint, e.g., in an LZ-End-like parsing, each phrase \( q_i \) is either (1) a symbol that is the leftmost occurrence of the symbol or (2) a (not necessary longest) prefix of the remaining suffix which is a suffix of \( q_1 \cdots q_j \) for some \( j < i \). We refer to an LZ-End-like parsing with the smallest number \( z_{end} \) of phrases, an optimal LZ-End parsing [12], and call the original, the greedy LZ-End parsing.\(^1\) Thus \( z \leq z_{no} \leq z_{end} \leq z_e \) holds.

Interestingly, \( z_{end} \leq g \) holds, where \( g \) is the size of the smallest context free grammar that derives (only) the string, while a similar relation between \( z_e \) and \( g \) does not seem to be known [12].

This brings us to two natural and important questions about the measure \( z_{end} \):

1. How efficiently can we compute \( z_{end} \)?
2. How much smaller can \( z_{end} \) be compared to \( z_e \)?

In this work, we answer a part of the above questions. Namely:

1. We prove the NP-hardness of computing \( z_{end} \).
2. We present an algorithm for exact computation by MAX-SAT.
3. We give a lower bound of the maximum value of the ratio \( z_e/z_{end} \).

In Section 3, we give the hardness result. Our reduction is from the vertex cover problem: finding a minimum set \( U \) of vertices such that every edge in a graph is incident to some vertex in \( U \). In Section 4, we show a MAX-SAT formulation for computing the optimal\(^1\) Notice that we do not need the distinction for LZ77, since the greedy LZ77 parsing is also an optimal LZ77-like parsing.
LZ-End parsing that follows an approach by Bannai et al. that allows computing NP-hard repetitiveness measures using MAX-SAT solvers [1]. In Section 5, we consider the ratio $z_e/z_{\text{end}}$. We give a family of binary strings such that the ratio asymptotically approaches 2. Note that we can easily modify this result to a larger alphabet. Since $(z_e/z_{\text{end}}) \leq (z_e/z_{\text{end}})$, the bound is tight, assuming that the conjecture by Kreft and Navarro [10] holds.

## Related work

The LZ77 and LZ78 are original members of the LZ family [15, 16]. It is well-known that the (greedy) LZ77 parsing produces the optimal version of the parsing [11]. The LZ78 parsing satisfies that each phrase can be represented as a concatenation of a previous phrase and a symbol. The NP-hardness of computing the optimal version of the LZ78 variant was shown [2]. This hardness result is also given by a reduction from the vertex cover problem. However, our construction of the reduction for the LZ-End differs from that for the LZ78 since these parsings have very different structures. The smallest string attractor [7] is one of the most fundamental repetitiveness measures. It is also known that computing the smallest string attractor of a given string is NP-hard [7]. The hardness result was proven by a reduction from the set-cover problem.

## 2 Preliminaries

### 2.1 Strings

Let $\Sigma$ be an alphabet. An element of $\Sigma^*$ is called a string. The length of a string $w$ is denoted by $|w|$. The empty string $\varepsilon$ is the string of length 0. Let $\Sigma^+ = \Sigma^* \setminus \{\varepsilon\}$. For any strings $x$ and $y$, $x \cdot y$ denotes the concatenation of two strings. We will sometimes abbreviate “-” (i.e., $x \cdot y = xy$). For a string $w = xyz$, $x$, $y$ and $z$ are called a prefix, substring, and suffix of $w$, respectively. They are called a proper prefix, a proper substring, and a proper suffix of $w$ if $x \neq w$, $y \neq w$, and $z \neq w$, respectively. The $i$-th symbol of a string $w$ is denoted by $w[i]$, where $1 \leq i \leq |w|$. For a string $w$ and two integers $1 \leq i \leq j \leq |w|$, let $w[i..j]$ denote the substring of $w$ that begins at position $i$ and ends at position $j$. For convenience, let $w[i..j] = \varepsilon$ when $i > j$. We will sometimes use $w[i..j]$ to denote $w[i..j-1]$. For any string $w$, let $w^1 = w$ and let $w^k = w w^{k-1}$ for any integer $k \geq 2$, i.e., $w^k$ is the $k$-times repetition of $w$.

### 2.2 LZ-End parsing

We give a definition of the LZ-End parsing, which is a variant of the Lempel-Ziv family.

**Definition 1** (Greedy LZ-End parsing). The LZ-End parsing of a string $w$ is the parsing $\text{LZEnd}(w) = q_1, \ldots, q_{z_e}$ of $w$ such that $q_i$ is either a symbol that is the leftmost occurrence of the symbol or the longest prefix of $q_i \cdots q_{z_e}$ that occurs as a suffix of $q_1 \cdots q_j$ for some $j < i$, which we call a source of the phrase.

We refer to each $q_i$ as a phrase. This definition, used in [8], is slightly different from the original version [9, 10] where a symbol is added to each phrase. The results in this paper hold for the original version as well (which we will show in the full version of the paper), but here we use this definition for simplicity. In this paper, we consider a more general version of the LZ-End parsing: a parsing $q_1, \ldots, q_{z_{\text{end}}}$ of a string $w$ such that $q_i$ is a (not necessary longest) suffix of $q_1 \cdots q_j$ for some $j < i$. We call such a parsing with a minimum number $z_{\text{end}}$ of phrases an optimal LZ-End parsing of $w$. We give an example of the greedy LZ-End parsing and the optimal LZ-End parsing in Figure 1.
Let $w = aacbbbbaababbabba$. The greedy LZ-End parsing $LZEnd(w)$ of $w$ is illustrated in the upper part of the figure. For the phrase at position 10, a longer substring $w[10..11] = ba$ has another previous occurrence at position 7, but there is no phrase that ends at position 8, and any longer substring does not have a previous occurrence. Therefore, the phrase starting at position 10 is $b$. The lower part of the figure shows an optimal LZ-End parsing (which is smaller than the greedy one) on the same string. Each phrase has a previous occurrence that ends at the end of some LZ-End phrase. The size of the greedy parsing is 12 and the size of the optimal parsing is 11.

### 2.3 Graphs

Let $G = (V, E)$ be a graph with the set of vertices $V$ and the set of edges $E$. An edge $e = \{u, v\}$ is called an incident edge of $u$. We denote the set of incident edges of $v$ as $\Gamma_G(v)$ and drop the subscript whenever it is clear from context. For an edge $e = \{u, v\}$, vertices $u$ and $v$ are the end points of $e$. For a subset of vertices $U \subseteq V$, $U$ is a vertex cover if for any $e \in E$, at least one end point of $e$ is contained in $U$. Let $\tau_G$ be the size of the minimum vertex cover of $G$ (i.e., $\tau_G$ denotes the vertex cover number of $G$). Notice that computing $\tau_G$ is NP-complete [3].

### 2.4 Maximum Satisfiability (MAX-SAT) problem

Let $\{x_1, \ldots, x_n\}$ be a set of literals and $C$ be a conjunctive normal form (CNF) formula. Each variable in $C$ is assigned a Boolean value (i.e., true or false). The goal of the Satisfiability (SAT) problem is to compute an assignment of variables that satisfies all clauses of $C$. The Maximum Satisfiability (MAX-SAT) problem is a variant of SAT, in which there are two types of clauses: hard clauses and soft clauses. A solution for MAX-SAT is a truth assignment of the variables such that all hard clauses are satisfied, and the number of soft clauses that are satisfied is maximized.

### 3 NP-hardness of computing the optimal LZ-End parsing

In this section, we consider the problem of computing the optimal LZ-End parsing of a given string. A decision version of the problem is given as follows.

- **Problem 2** (Decision version of computing the optimal LZ-End parsing (OptLE)). Given a string $w$ and an integer $k$, decide whether there exists an LZ-End parsing of size $k$ or less.

We show the NP-completeness of OptLE in the following and present an algorithm for exact computation in the next section.

- **Theorem 3.** OptLE is NP-complete.

**Proof.** We give a reduction from the vertex cover problem to OptLE. Let $G = (V, E)$ be a graph with a set of vertices $V = \{v_1, \ldots, v_n\}$ and a set of edges $E = \{e_1, \ldots, e_m\}$. Suppose that an input graph $G$ of the vertex cover problem is connected and $|\Gamma(v)| \geq 2$ for any
Let $G = (V, E)$ be the complete graph of three vertices $v_1, v_2, v_3$ and $e_1 = \{v_1, v_2\}$, $e_2 = \{v_2, v_3\}$, $e_3 = \{v_1, v_3\}$. $W_G$ and the greedy parsing and an optimal parsing are illustrated in the figure. The first two parts ($P$ and $Q$) share the same parsing. The last two parts ($R$ and $S$) are different. The upper part in the figure shows the greedy parsing and the lower part shows an optimal parsing. For instance, in the optimal parsing, we can choose $e_1^3$ and $e_2^3$ as phrases by using non-greedy parsing in $R_1$. In other words, we can reduce two phrases in $S$-part by adding one phrase in $R_1$. In this example, the optimal parsing represents a vertex cover $\{v_1, v_3\} \subset V$ of $G$ (since $R_2$ selects the greedy parsing and the others are not).

$v \in V$. We identify each vertex $v_i$ as a symbol $v_i$ and each edge $e_j$ as a symbol $e_j$. We also introduce the symbol $\$, and a set of symbols that occur uniquely in the string. The latter is represented, for simplicity, by the special symbol $\#$, i.e., $\#$ represents a different symbol each time it occurs in our description. We consider the string $W_G$ defined by graph $G$ as follows.

- $W_G = \prod_{i=1}^{m} P_i \cdot \prod_{i=1}^{m} Q_i \cdot \prod_{i=1}^{m} R_i \cdot \prod_{i=1}^{m} S_j$
- $P_i = e_1^3 \# v_1^2 \# v_2^2 \# v_3^2 \# \chi_i \cdot \gamma_i$
- $Q_i = e_1^3 \#$
- $R_i = e_1^3 \prod_{e_j \in \Gamma(v_i)} (e_j^3 v_j^2 \#)$
- $S_j = e_2^3 \#$
- $\chi_i = \prod_{e_j \in \Gamma(v_i)} (v_i e_j \#)$
- $\gamma_i = \prod_{e_j \in \Gamma(v_i)} (e_j^2 v_j \#)$

An example of this string is illustrated in Figure 2. Note that we use $i$ for representing indices of vertices and $j$ for indices of edges.

Before we show the detail of the reduction, we start with an intuitive description of our reduction. The string $W_G$ consists of four parts (which are represented by $P$, $Q$, $R$, and $S$).

1. The first two parts are non-functional parts. They can only be parsed in a single sensible way. These parts play a role as sources of the third part ($R$-part).
2. In the third part, $R_i$ corresponds to the vertex $v_i$. Roughly speaking, $R_i$ can be parsed in two sensible ways such that the parsing represents whether the vertex $v_i$ is in the vertex cover or not. If a vertex $v_i$ is in the vertex cover, then the parsing of $R_i$ needs one more phrase.
3. In the last part, $S_j$ corresponds to the edge $e_j$. $S_j$ will be parsed into two phrases iff one of the incident vertex of the edge $e_j$ is in the vertex cover. Otherwise, $S_j$ has three phrases. Overall, minimizing the number of vertices for the vertex cover of a graph $G$ corresponds to minimizing the total penalties in the $R$-part such that every $S_j$ will be parsed into two phrases.
We show that the number of phrases of the optimal parsing of $W_G$ is less than $13n + 22m + k$ if and only if the vertex cover number $\tau_G$ is less than $k$.

First, we observe an optimal LZ-End parsing of $W_G$. Let us consider a parsing of $\prod_{i=1}^n P_i$. In this part, the greedy parsing gives $10n + 13m$ phrases. In the greedy parsing of $\prod_{i=1}^n P_i$, phrases $v_i^2$ in $v_i^2 \#$, $v_i$ in $v_i^2 \#$, and the second occurrence of $v_i^2$ have length 2, and the other phrases have length 1. It is easy to see that this parsing is a smallest possible parsing of $\prod_{i=1}^n P_i$. Moreover, other parsings of the same size do not affect the parsing of the rest of the string; candidates for a source cannot be increased by selecting any other parsings since the phrases of length 2 are preceded by unique symbols #. Hence, we can choose this greedy parsing as a part of an optimal parsing.

In the second part $\prod_{j=1}^m Q_j$, the greedy parsing also gives an optimal parsing which has $3m$ phrases (i.e., each $Q_j$ is parsed into three phrases since $e_j^3$ occurs in $P_i$ for some $i$ and $e_j^3$ is unique in $\prod_{j=1}^m Q_j$). This parsing is also a smallest possible parsing and does not affect any parsings of the rest of the string.

The remaining suffix $\prod_{j=1}^m R_j \cdot \prod_{j=1}^m S_j$ is a key of the reduction. The key idea is that $S_j$ represents whether the edge $e_j$ is an incident edge of some vertex in a subset of vertices or not. $\#e_j^3$ in $S_j$ has exactly two previous occurrences in the $R$-part (since each edge is incident to exactly two vertices). Hence $S_j$ can be parsed into two phrases (i.e., $\#e_j^3$, #) if and only if $\#e_j^3$ has an occurrence which ends with an LZ-End phrase in the $R$-part. Now we consider the greedy parsing of the $R_j$-part (let $\Gamma(v_i) = \{e_{(i,1)}, \ldots, e_{(i,|\Gamma(v_i)|)}\}$), which is as follows:

$$v_i^3, \ldots, v_i^3, e_{(i,1)}, \ldots, v_i^3, e_{(i,|\Gamma(v_i)|)} e_{(i,|\Gamma(v_i)|)}^2 v_i, v_i^2, \#.$$  

The parsing has $2|\Gamma(v_i)| + 3$ phrases. We claim that this parsing is the smallest possible parsing: If the length of every phrase is at most 3, then $2|\Gamma(v_i)| + 3$ is the minimum size since the length of $R_i$ is $6|\Gamma(v_i)| + 7$. On the other hand, we can see that substrings of length at least 4 which contain a symbol $v_i$ are unique in the whole string $W_G$ by the definition. Namely, $\#e_j^3$ is the only substring of length at least 4 which is not unique. Let us consider a parsing of $R_i$ such that the parsing has $\alpha$ length-4 phrases. In other words, we choose $\alpha$ incident edges out of $|\Gamma(v_i)|$ edges. Let $(i_1, \ldots, i_\alpha)$ be the sequence of indexes of selected edges. We observe that the length of substrings that are covered by length at most 3 phrases. The length of the prefix of $R_i$ that is succeeded by the first length-4 phrase $\#e_{(i_1,i_1)}^3$ is $6(i_1 - 1) + 4$. This implies that there are at least $2(i_1 - 1) + 2$ phrases. The length of substring between $\#e_{(i_d,i_d)}$ and $\#e_{(i_{d+1},i_{d+1})}$ is $6(i_d - i_{d-1}) + 2$. Thus there are at least $2(i_d - i_{d-1}) + 1$ phrases in each middle part. The length of the suffix that is preceded by the last length-4 phrase $\#e_{(i_\alpha,i_\alpha)}^3$ is $6(|\Gamma(v_i)| - i_\alpha) + 5$. Since the last symbol is a unique symbol #, there are at least $2(|\Gamma(v_i)| - i_\alpha) + 3$ phrases in the last part. Hence, there are at least

$$\alpha + 2(i_1 - 1) + 2 + \sum_{d=2}^\alpha (2(i_d - i_{d-1}) + 1) + 2(|\Gamma(v_i)| - i_\alpha) + 3 = 2|\Gamma(v_i)| + 4$$

phrases. Thus the minimum number of phrases of $R_i$ is $2|\Gamma(v_i)| + 3$ and the above greedy parsing is the only candidate which is the minimum size. Notice that phrases of this parsing do not end with $\#e_j^3$. Let us consider the other possible parsing of $R_i$-part as follows:

$$v_i^2, v_i^2, e_{(i,1)}, \ldots, v_i^2, e_{(i,|\Gamma(v_i)|)} e_{(i,|\Gamma(v_i)|)}^2, v_i^2, \#, \#.$$  

This parsing has $2|\Gamma(v_i)| + 4$ phrases. Notice that this parsing has phrases which end with $\#e_j^3$.

Thus $S_j$ can be parsed into two phrases if we choose a non-greedy parsing such that there exists a phrase that ends at one of these positions. In other words, if we choose such a parsing
in the \( R_i \)-part, we can reduce at most \( |\Gamma(v_i)| \) phrases in the \( S \)-part. These observations imply that \( R_i \) is parsed into \( 2|\Gamma(v_i)| + 3 \) or \( 2|\Gamma(v_i)| + 4 \) phrases in any optimal parsing of \( W_G \).

Let us consider an optimal LZ-End parsing. Let \( r \) be the number of substrings \( R_i \) which contain \( 2|\Gamma(v_i)| + 4 \) phrases, and \( s \) be the number of substrings \( S_j \) which contain exactly two phrases. Then the size of the parsing is

\[
(10n + 13m) + (3m) + (2 \sum_{i=1}^{n} |\Gamma(v_i)| + 3n + r) + (3m - s) = 13n + 23m + r - s.
\]

We consider a subset \( V' \) of vertices such that \( v_i \in V' \) if and only if \( R_i \) is parsed into \( 2|\Gamma(v_i)| + 4 \) phrases (i.e., \( |V'| = r \)), and a subset \( E' \) of edges such that \( e_j \in E' \) if and only if \( S_j \) is parsed into three phrases (i.e., \( |E'| = m - s \)). If \( E' = \emptyset \) (i.e., \( s = m \)), \( V' \) is a vertex cover of \( G \). Otherwise, \( V' \) is not a vertex cover of \( G \). However we can obtain the vertex cover number by using the parsing. Since the parsing is an optimal parsing, we can observe that there is no vertex \( v_i \) in \( V \setminus V' \) which has two or more incident edges in \( E' \) (we can reduce two or more phrases in \( S \)-part by adding one phrase in \( R_i \), a contradiction). This implies that we can obtain a vertex cover by choosing one vertex in \( V \setminus V' \) for each edge in \( E \setminus E' \). Then there exists an optimal LZ-End parsing of the same size which can directly represent a vertex cover. In other words, the vertex cover number is \( r + m - s \) if there exists an optimal LZ-End parsing of \( 13n + 22m + (r + m - s) \) phrases. It is clear from the above constructions that there exists an optimal LZ-End parsing of \( 13n + 22m + k \) phrases iff the vertex cover number is \( k \).

Since we can check a parsing is an LZ-End parsing in linear time, \( \text{OptLE} \) is clearly in class NP.

### 4 MAX-SAT Formulation

An approach for exact computation of various NP-hard repetitiveness measures was shown in [1], where they formulated them as MAX-SAT instances so that very efficient solvers could be taken advantage of. Here, we show that this approach can be adapted to computing the optimal LZ-End parsing as well.

Let the input string be \( T[1..n] \), and for any \( i \in [2, n] \), let \( M_i = \{ j \mid 1 \leq j < i, T[j] = T[i] \} \).

Below, we use 1 to denote true, and 0 to denote false. We introduce the following Boolean variables:

- \( p_i \) for all \( i \in [1, n] \): \( p_i = 1 \) if and only if position \( i \) is a starting position of an LZ-End phrase. Note that \( p_1 = 1 \).
- \( c_i \) for all \( i \in [1, n] \): \( c_i = 1 \) if and only if position \( i \) is the left-most occurrence of symbol \( T[i] \).
- \( r_{i \rightarrow j} \) for all \( i \in [2, n] \) and \( j \in M_i \): \( r_{i \rightarrow j} = 1 \) if and only if position \( i \) references position \( j \) via an LZ-End factor.

Notice that the truth values of \( c_i \) are all fixed for a given string and are easy to determine. Furthermore, the left-most occurrence must be beginning of a phrase, so, some values of \( p_i \) can also be fixed. For all \( i \in [1, n] \):

\[
\begin{align*}
c_i &= p_i = 1 \text{ if } i \text{ is left-most occurrence of } T[i], \\
c_i &= 0 \text{ otherwise.}
\end{align*}
\]
The truth values of \( p_i \) define the factors, so in order to minimize the number of factors, we define the soft clauses as \( \overline{p_i} \) for all \( i \in [1, n] \). Below, we give other constraints between the variables that must be satisfied, i.e., hard clauses.

The symbol at any position must either be a left-most occurrence, or it must reference some position to its left. That is, for any \( i \in [1, n] \):

\[
e_i + \sum_{j \in M_i} r_{i \rightarrow j} = 1. \tag{3}
\]

In order to ensure that references in the same LZ-End phrase are consistent, we have the following two constraints. The first ensures that if \( i \) references \( j \) and the symbols at positions \( i - 1 \) and \( j - 1 \) are different or do not exist (i.e., \( j = 1 \)), positions \( i \) and \( i - 1 \) cannot be in the same LZ-End phrase. For all \( i \in [2, n] \) and \( j \in M_i \) such that \( j = 1 \) or \( T[j - 1] \neq T[i - 1] \):

\[
r_{i \rightarrow j} \implies p_i, \tag{4}
\]

The second ensures that if position \( i \) references position \( j \) and \( i \) is not a start of an LZ-End phrase, then, position \( i - 1 \) must reference position \( j - 1 \). For all \( i \in [2, n] \) and \( j \in M_i \) \( \setminus \{1\} \) s.t. \( T[j - 1] = T[i - 1] \):

\[
r_{i \rightarrow j} \land \overline{p_i} \implies r_{i - 1 \rightarrow j - 1}. \tag{5}
\]

Finally, the following constraints ensure that the reference of each LZ-End phrase must end at an end of a previous LZ-End phrase. For all \( i \in [1, n] \) and \( j \in M_i \):

\[
\begin{cases}
r_{i \rightarrow j} \land p_{i + 1} \implies p_{j + 1} & \text{if } i \in [1, n) \\
r_{i \rightarrow j} \implies p_{j + 1} & \text{if } i = n. \tag{6}
\end{cases}
\]

It is easy to see that the truth assignments that are derived from any LZ-End parsing will satisfy the above constraints.

We now show that any truth assignment that satisfies the above constraints will represent a valid LZ-End parsing. The truth values for \( p_i \) implies a parsing where each phrase starts at a position \( i \) if and only if \( p_i = 1 \). Constraint (1),(2),(3) ensure that each position is either a left-most occurrence or references a unique previous position. Thus, it remains to show that the referencing of each position of a given factor is consistent (adjacent positions reference adjacent positions) and ends at a previous phrase end.

For any position \( i \) such that \( e_i = 0 \), let \( j \in M_i \) be the unique value such that \( r_{i \rightarrow j} = 1 \). We can see that any such position \( i \) that is not at the beginning of a phrase (i.e., \( p_i = 0 \)) will reference a position consistent with the reference of position \( i - 1 \): If \( j = 1 \) or \( T[j - 1] \neq T[i - 1] \), then Constraint (4) would imply \( r_{i \rightarrow j} = 0 \). Thus, we have \( j > 1 \) and \( T[j - 1] = T[i - 1] \), and from Constraint (5), we have that \( r_{i - 1 \rightarrow j - 1} \), and the referencing inside a factor is consistent. Finally, from Constraint (6), the last reference in a phrase always points to an end of a previous LZ-End phrase.

The MAX-SAT instance contains \( O(n^3) \) variables, and the total size of the CNF is \( O(n^2) \): \( O(n) \) clauses of \( O(n) \) size (Constraint (3) using linear size encodings of cardinality constraints, e.g. [14]), and \( O(n^3) \) clauses of size \( O(1) \) (the soft clauses, and Constraints (4), (5), (6)).

We note that it is not difficult to obtain a MAX-SAT formulation for the original definition of LZ-End by minor modifications.
5 Approximation ratio of greedy parsing to optimal parsing

In this section, we consider an approximation ratio of the size $z_e$ of the greedy LZ-End parsing to the size $z_{end}$ of the optimal LZ-End parsing. Here, we give a lower bound of the ratio.

Theorem 4. There exists a family of binary strings such that the ratio $z_e/z_{end}$ asymptotically approaches 2.

Proof. Let $K = \sum_{i=1}^{k} 2^i (\leq 2^{k+1} - 2)$ for any positive integer $k \geq 1$. The following binary string $w_k$ over an alphabet $\{a, b\}$ gives the lower bound:

$$w_k = aa \cdot \prod_{i=1}^{k} (a^{2^i}) \cdot b \cdot \prod_{i=1}^{K} (a^i b^3).$$

It is easy to see that $K$ is the length of the substring $\prod_{i=1}^{k} (a^{2^i})$. First, we show the greedy parsing of $w_k$. Let $W_0 = aa \cdot \prod_{i=1}^{k} (a^{2^i}) \cdot b^4$ (i.e., a prefix of $w_k$) and $W_j = W_{j-1} \cdot a^j b^3$ for any $1 \leq j \leq K$. Notice that $W_k = w_k$. We show that

$$LZEnd(W_j) = LZEnd(W_{j-1}) \cdot a^j b^3, b$$

(7)

by induction on $j$. Initially, we consider the greedy parsing of $W_0$. The greedy parsing of the first run (i.e., maximal substring with a unique symbol) is $a, a, a^2, \ldots, a^{2^k}$ of size $k + 2$. The second run is parsed into three phrases $b, b, b^2$. Thus

$$LZEnd(W_0) = a, a, a^2, \ldots, a^{2^k}, b, b, b^2.$$  

Moreover, we can see that the greedy parsing of $W_1$ is

$$LZEnd(W_1) = a, a, a^2, \ldots, a^{2^k}, b, b, b^2, ab^2, b.$$  

Hence Equation 7 holds for $j = 1$. Suppose that Equation 7 holds for any $j \leq p$ for some integer $p \geq 1$. We show that Equation 7 holds for $j = p + 1$. Assume that there exists a phrase $x$ of $LZEnd(W_{p+1})$ which begins in $W_p$ and ends in a new suffix $a^{p+1} b^3$ of $W_{p+1}$. By the induction hypothesis, phrases of $LZEnd(W_p)$ which end with $a$ are only in the first $a$’s run. This implies that $x$ cannot end with $a$ and $x$ can be written as $x = x' b a^{p+1} b^\ell$ for some prefix $x'$ of $x$ and some positive integer $\ell$. However, $a^{p+1}$ only occurs in the first $a$’s run. Thus $LZEnd(W_{p+1})$ cannot have such a phrase $x$, namely $LZEnd(W_{p+1}) = LZEnd(W_p)$, $S$ for some factorization $S$ of the remaining suffix $a^{p+1} b^3$. It is easy to see that the remaining suffix $a^{p+1} b^3$ of $W_{p+1}$ is parsed into $a^{p+1} b^2, b$. Hence Equation 7 holds for $j = p + 1$, and it also holds for any $j$. Notice that $|LZEnd(w_k)| = 2K + k + 5$ holds.

Finally, we give a smaller parsing of $w_k$. We consider the same parsing for the first run and a different parsing for the second run as $b, b, b, b$. In the greedy parsing, $a^i b^3$ cannot be a phrase since the only previous occurrence does not have an LZ-End phrase. We can use a substring $a^i b^3$ as a new phrase of $W_j$ (see also Figure 3). Thus there exists an LZ-End parsing

$$a, a, a^2, \ldots, a^{2^k}, b, b, b, b, a^i b^3, \ldots, a^K b^3.$$  

The size of the parsing is $K + k + 6$. Therefore the ratio $z_e/z_{end}$ asymptotically approaches 2 for this family of strings.

Note that this family of strings also gives a lower bound of the ratio $z_e/z_{na}$ since $(z_e/z_{end}) \leq (z_e/z_{na})$ holds.
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Figure 3 Illustration for two variants of LZ-End parsings of a string $w_k$ (Theorem 4). In the optimal parsing, we can choose $a^j b^3$ (dotted lines) as a phrase for each $j$ ($1 \leq j \leq K$) by adding a single letter phrase $b$.

6 Conclusions

In this paper, we first studied the optimal version of the LZ-End variant. We showed the NP-completeness of the decision version of computing the optimal LZ-End parsing and presented an approach for exact computation of the optimal LZ-End by formulating as MAX-SAT instances. We also gave a lower bound of the possible gap (as the ratio) between the greedy LZ-End and the optimal LZ-End. Finally, we note possible future work in the following.

- Our reduction from the vertex cover problem uses a polynomially large alphabet. How can we construct a reduction with a small alphabet?
- The most interesting remaining problem is an upper bound of the ratio discussed in Section 5. We conjecture that there exists a constant upper bound (i.e., $z_e/z_{end} \leq c$ for any strings where $c$ is a constant). This implies that the greedy parsing gives a constant-approximation of the optimal parsing. On the other hand, if there exists a family of strings which gives $c > 2$ or non-constant ratio, then the conjecture $z_e \leq 2z_{no}$ does not stand.

References


Sliding Window String Indexing in Streams

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Abstract

Given a string $S$ over an alphabet $\Sigma$, the string indexing problem is to preprocess $S$ to subsequently support efficient pattern matching queries, that is, given a pattern string $P$ report all the occurrences of $P$ in $S$. In this paper we study the streaming sliding window string indexing problem. Here the string $S$ arrives as a stream, one character at a time, and the goal is to maintain an index of the last $w$ characters, called the window, for a specified parameter $w$. At any point in time a pattern matching query for a pattern $P$ may arrive, also streamed one character at a time, and all occurrences of $P$ within the current window must be returned. The streaming sliding window string indexing problem naturally captures scenarios where we want to index the most recent data (i.e. the window) of a stream while supporting efficient pattern matching.

Our main result is a simple $O(w)$ space data structure that uses $O(\log w)$ time with high probability to process each character from both the input string $S$ and any pattern string $P$. Reporting each occurrence of $P$ uses additional constant time per reported occurrence. Compared to previous work in similar scenarios this result is the first to achieve an efficient worst-case time per character from the input stream with high probability. We also consider a delayed variant of the problem, where a query may be answered at any point within the next $\delta$ characters that arrive from either stream. We present an $O(w + \delta)$ space data structure for this problem that improves the above time bounds to $O(\log(w/\delta))$. In particular, for a delay of $\delta = \epsilon w$ we obtain an $O(w)$ space data structure with constant time processing per character. The key idea to achieve our result is a novel and simple hierarchical structure of suffix trees of independent interest, inspired by the classic log-structured merge trees.

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Introduction

The string indexing problem is to preprocess a string $S$ into a compact data structure that supports efficient subsequent pattern matching queries, that is, given a pattern string $P$, report all occurrences of $P$ within $S$. In this paper, we introduce a basic variant of string indexing called the streaming sliding window string indexing (SSWSI) problem. Here, the string $S$ arrives as a stream one character at a time, and the goal is to maintain an index of a window of the last $w$ characters, for a specified parameter $w$. At any point in time a pattern matching query for a pattern $P$ may arrive, also streamed one character at a time, and we need to report the occurrences of $P$ within the current window. The goal is to compactly maintain the index while processing the characters arriving in either stream efficiently. We consider two variants of the problem: a timely variant where each query must be answered immediately, and a delayed variant where it may be answered at any point within the next $\delta$ characters arriving from either stream, for a specified parameter $\delta$. See Section 1.1 for precise definitions.

The SSWSI problem naturally captures scenarios where we want to index the most recent data (i.e. the window) of a stream while supporting efficient pattern matching. For instance, monitoring a high-rate data stream system where we cannot feasibly index the entire stream but still want to support efficient queries. Depending on the specific system we may require immediate answers to queries, or we may be able to afford a delay that allows for more efficient queries and updates.

The SSWSI problem has not been explicitly studied before in our precise formulation, but for the timely variant several closely related problem are well-studied. In particular, the sliding window suffix tree problem \cite{8,13,24,25,28} is to maintain the suffix tree of the current window (i.e., the compact trie of the suffixes of the window) as each character arrives. With appropriate augmentation the suffix tree can be used to process pattern matching queries efficiently, leading to a solution to the timely SSWSI problem. For constant-sized alphabets, the best of these solutions \cite{8} maintains the sliding window suffix tree in constant amortized time per character while supporting efficient pattern matching queries. The worst-case time for updates is $\Omega(w)$. The other solutions achieve similar amortized time bounds. This amortization cannot be avoided since explicitly maintaining the suffix tree after the arrival of a new character may incur $\Omega(w)$ changes.

Another closely related problem is the online string indexing problem \cite{3,4,5,7,14,21,22,23}. Here the goal is to process $S$ one character at a time (in either left-to-right or right-to-left order), while incrementally building an index of the string read so far. The best of these solutions update the index in either constant time per character for constant-sized alphabets \cite{23} or $O(\log \log n + \log \log |\Sigma|)$ time for any alphabet where each character fits in a constant number of machine words \cite{21}. These solutions all heavily rely on processing the string in right-to-left order to avoid the inherent linear time suffix tree updates due to appending, as mentioned above. Therefore they cannot be applied in our left-to-right streaming setting. Alternatively, we can instead apply these solutions on the reverse of the string $S$, but then each pattern must be processed in reverse order, which also cannot be done in our setting. Also, note that these solutions index the entire string read so far. It is not clear if they can be adapted to efficiently index a sliding window.

Another line of work shows how to maintain a fully dynamic suffix array under insertions and deletions \cite{1,2,19,27}. These can also be used to solve SSWSI but are more general and lead to polylogarithmically slower bounds than our results while being more complicated.
Our main result is an efficient and simple solution to the SSWSI problem in both the timely and delayed variant. Let \( w \) denote the size of the window. For the timely variant, we present a string index that uses \( O(w) \) space and processes a character from the stream \( S \) in \( O(\log w) \) time. Each pattern matching query \( P \) is also supported in \( O(\log w) \) time per character with additional \( O(\text{occ}) \) time incurred after receiving the last character of \( P \), where \( \text{occ} \) is the number of occurrences of \( P \) in the current window. The index is randomized and both time bounds hold with high probability. Compared to previous suffix tree based approaches for indexing a sliding window, we improve the worst-case time bounds per character in the stream from \( \Omega(w) \) to \( O(\log w) \) with high probability. This is particularly important in the above mentioned applications, such as high-rate data stream systems. Our solution generalizes to the delayed variant of the problem. If we allow a delay of \( \delta \) before answering each query we achieve \( O(w + \delta) \) space while improving the above time bounds to \( O(\log(w/\delta)) \). In particular, if we allow a delay of \( \delta = \epsilon w \) for any constant \( \epsilon > 0 \), we achieve linear space and optimal constant time (reporting the occurrences still takes \( O(\text{occ}) \) time, and we do not count the reporting time towards the delay). Note that \( \delta \leq w \) is sufficient delay for optimal time bounds and we can assume \( O(w + \delta) = O(w) \). The results hold on a word RAM and for any alphabet size, assuming that each character fits into a constant number of machine words.

The key idea to achieve our result is a novel and simple hierarchical structure of suffix trees inspired by log-structured merge trees [26]. Instead of maintaining a single suffix tree on the window we maintain a collection of suffix trees of exponentially increasing sizes that cover the current window. We show how to efficiently maintain the structure as new characters from the stream arrive by incrementally “merging” suffix trees, while supporting efficient pattern matching queries within the window.

Our solution uses randomization to construct suffix trees in linear time with high probability. Plugging in a deterministic construction algorithm such as the one by Ukkonen [30], we obtain a solution using \( O(\log w \log |\Sigma|) \) time for both queries and updates. With more recent deterministic suffix tree solutions [6, 10, 14] we can improve this to obtain \( O(\log w \log \log n) \) time per character for both queries and updates. Note that the \( O(\log \log |\Sigma|) \) in the time bounds of [14] has been replaced by \( O(\log \log n) \) here due to an additional sorting step using [17].

1.1 Setup and Results

We formally define the problem as follows. Let \( S \) be a stream over any alphabet \( \Sigma \) where each character fits in a constant number of machine words. For given integer parameters \( w \geq 1 \) and \( \delta \geq 0 \), the \( \delta \)-delayed streaming sliding window string indexing ((w, \delta)-SSWSI) problem is to maintain a data structure that, after receiving the first \( i \) characters of \( S \), supports:

- **Report(P)**: report all the occurrences of \( P \) in \( S[i - w + 1, i] \) before an additional \( \delta \) characters have arrived from either stream.
- **Update()**: process the next character in the stream \( S \).

In the Report(P) query the pattern string \( P \) is also streamed. When \( P \) is streamed it interrupts the stream \( S \), arrives one character at a time, and all characters of \( P \) arrive before the streaming of \( S \) resumes. Furthermore, we do not assume that we know the length of \( P \) before the arrival of its last character. Although \( P \) is streamed we assume random access to its characters after they arrive, as any pattern that fits in the window is at most \( w \) characters long and we can afford to store it. The delay is counted from after the last character of \( P \) arrives. Characters from \( S \) and from new patterns count towards the delay, while reported occurrences do not (otherwise it would be impossible to answer the query in time if there are more than \( \delta \) occurrences).
We define the *timely streaming sliding window string indexing* (w-SSWSI) problem to be \((w, 0)-SSWSI\), that is, queries must be answered immediately as the last character of the pattern arrives.

We show the following general main result.

\[\text{Theorem 1.}\] Let \(S\) be a stream and let \(w \geq 1\) and \(\delta \geq 0\) be integers. We can solve the \((w, \delta)-SSWSI\) problem on \(S\) with an \(O(w + \delta)\) space data structure that supports \text{Update} and \text{Report} in \(O(\log \frac{w}{\delta} + 1)\) time per character with high probability. Furthermore, \text{Report} uses additional worst-case constant time per reported occurrence.

Here, with high probability means with probability at least \(1 - \frac{1}{w^d}\) for any constant \(d\).

Theorem 1 provides a trade-off in the delay parameter \(\delta\). In particular, plugging in \(\delta = 0\) in Theorem 1 we obtain a solution to the timely SSWSI problem that uses \(O(w)\) space and \(O(\log w)\) time per character for both \text{Update} and \text{Report}. Compared to the previous work on sliding window stream indexing \([8, 13, 18, 24, 25, 28, 29]\) this improves the worst-case bounds on the \text{Update} operation from \(\Omega(w)\) to \(O(\log w)\) with high probability and also removes the restriction on the alphabet. At the other extreme, plugging in \(\delta = \epsilon w\) for constant \(\epsilon > 0\) in Theorem 1 we obtain a solution to the delayed SSWSI problem that uses \(O(w)\) space and optimal constant time per character with high probability. All our results hold on a word RAM where each machine word has at least \(\log w\) bits, and where each character of the alphabet fits into a constant number of machine words.

### 1.2 Techniques

We obtain our result for the timely variant, but without high probability guarantees, as follows. At all times we maintain at most \(\log w\) suffix trees that do not overlap and together cover the window. The trees are organized by the log-structured merge technique \([26]\), where the rightmost tree is the smallest and their sizes increase exponentially towards the left. For each new character that arrives we append its suffix tree to the right side of our data structure. Whenever there are two trees of the same size next to each other we “merge” them by constructing a new suffix tree covering them both. Each character from \(S\) is involved in at most \(\log w\) merges and each merge takes expected linear time, so we spend expected amortized \(O(\log w)\) time per character in \(S\). We deamortize the updates by temporarily keeping both trees while merging them in the background. Note that for each adjacent pair of suffix trees we also store a suffix tree approximately covering them both, referred to as boundary trees (see details below).

We find the occurrences of a pattern \(P\) in the window by querying each of these trees, which takes \(O(\log w)\) time per character in \(P\). For adjacent pairs of trees larger than \(|P|\) we find the occurrences of \(P\) crossing from one into the other using the boundary trees. The remaining trees cover a suffix of the window of length \(O(|P|)\), and we grow a suffix tree to answer queries in this suffix at query time. Our data structure has some “overhang” on the left side of the window, and we use range maximum queries to report only the occurrences that start inside the window.

This solution is generalized to incorporate a delay of \(\delta\) as follows. We store the \(O(\log(\frac{w}{\delta}))\) largest trees from the timely solution and leave a suffix of size \(\Theta(\delta)\) of the window uncovered by suffix trees. We answer queries as follows. If \(|P| > \delta/4\) we say that \(P\) is long, and otherwise it is short. For long patterns we do as in the timely case; the suffix tree we grow at query time now must also contain the uncovered suffix, but it still has size \(O(|P|)\) since the uncovered part of the window has length \(O(\delta) = O(|P|)\). We show how to do this in \(O(\log(\frac{w}{\delta}))\) time per character in \(P\). For short patterns we utilize that they are smaller
than the delay to temporarily buffer the queries and later batch process them. We buffer up to $O(\delta \log(w/\delta))$ work and deamortize it over $\Theta(\delta)$ characters, obtaining the same bound as for long patterns. Updates run in the same bound since each character from $S$ is involved in at most $O(\log(w/\delta))$ merges before it leaves the window.

Finally, we improve the time bounds by proving that for any substring $S'$ of our window, we can construct the suffix tree over $|S'|$ time with probability $1 - w^{-d}$ for any constant $d > 1$. We do so by reducing the alphabet $\Sigma' = \{ c \in S' \}$ of $S'$ to rank-space $\{1, 2, \ldots, |\Sigma'|\}$ from which the algorithm by Farach-Colton et al. [12] can construct the suffix tree in worst-case linear time. For large strings ($|S'| > w^{1/5}$) we pick a hash function from $\Sigma \rightarrow [0, w]$ that with high probability is injective on $S'$, and then we use radix sort to reduce to rank-space in linear time. For small strings ($|S'| \leq w^{1/5}$) we pick a hash function from $\Sigma \rightarrow [0, w/\log w]$ that is injective with (almost) high probability, and use this to manually construct a mapping into rank space in $O(S')$ time. This mapping algorithm uses additional $O(w/\log w)$ space, but we construct at most $O(\log w)$ suffix trees at any time so the total space is linear.

1.3 Outline

In Section 2 we cover the preliminaries, including some useful facts about suffix trees. In Section 3 we give a solution to the timely SSWSI problem that supports each operation in expected logarithmic time per character. In Section 4 we show how to generalize this to incorporate delay, and in Section 5 we show how to get good probability guarantees, proving Theorem 1.

2 Preliminaries

Given a string $X$ of length $n$ over an alphabet $\Sigma$, the $i$th character is denoted $X[i]$ and the substring starting at $X[i]$ and ending at $X[j]$ is denoted $X[i, j]$. The substrings of the form $X[i, n]$ are the suffixes of $X$.

A segment of $X$ is an interval $[i, j] = \{i, i+1, \ldots, j\}$ for $1 \leq i \leq j \leq n$. We will sometimes refer to segments as strings, i.e., the segment $[i, j]$ refers to the string $X[i, j]$. The definition differs from “substring” by being specific about position; even if $X[1, 2] = X[3, 4]$ we have $[1, 2] \neq [3, 4]$. A segmentation of $X$ is a decomposition of $X$ into disjoint segments that cover it. For instance, $x_1 = [1, i]$ and $x_2 = [i+1, n]$ is a segmentation of $X$ into two parts. The two segments $x_1$ and $x_2$ are adjacent since $x_2$ starts immediately after $x_1$ ends, and for a pair of adjacent segments we define the boundary $(x_1, x_2)$ to be the implicit position between $i$ and $i+1$.

The suffix tree [31] $T$ over $X$ is the compact trie of all suffixes of $XS$, where $S \not\in \Sigma$ is lexicographically smaller than any letter in the alphabet. Each leaf corresponds to a suffix of $X$, and the leaves are ordered from left to right in lexicographically increasing order. The suffix tree uses $O(n)$ space by implicitly representing the string associated with each edge using two indices into $X$. Farach-Colton et al. [12] show that the optimal construction time for $T$ is $\operatorname{sort}(n, |\Sigma|)$, i.e., the time it takes to sort $n$ elements from the universe $\Sigma$. For alphabets of the form $\Sigma = \{0, \ldots, n\}$ for constant $c \geq 1$ this implies that $T$ can be built in worst-case $O(n)$ time using radix sort. For larger alphabets we can reduce to the polynomial case in expected linear time using hashing, building $T$ in expected linear time (see Section 5 for details).

The suffix array $L$ of $X$ is the array where $L[i]$ is the starting position of the $i$th lexicographically smallest suffix of $X$. Note that $L[i]$ corresponds to the $i$th leaf of $T$ in left-to-right order. Furthermore, let $v$ be an internal node in $T$ and let $s_v$ be the string
spelled out by the root-to-$v$ path. The descendant leaves of $v$ exactly correspond to the suffixes of $X$ that start with $s_v$, and these leaves correspond to a consecutive range $[\alpha, \beta]_v$ in $L$.

We augment the suffix tree to support efficient pattern matching queries as follows. First, we use the well-known FKS perfect hashing scheme [15] to store the edges of the suffix tree, so we can for any node determine if there is an outgoing edge matching a character $a \in \Sigma$ in worst-case constant time. Note that this construction takes expected linear time. Furthermore, we also build a range maximum query data structure over $L$. This data structure supports range maximum queries, i.e., given a range $[\alpha, \beta]$ return the $j \in [\alpha, \beta]$ maximizing $L[j]$. It also supports range minimum queries, defined analogously. The data structure can be built in linear time and supports queries in constant time [16]. Finally, we preprocess the suffix tree in linear time such that each internal node $v$ stores the range $[\alpha, \beta]_v$ into $L$ corresponding to the occurrences of $s_v$.

We can use this structure to efficiently find all the occurrences of $P$ in $O(|P| + \text{occ})$ time, where $\text{occ}$ is the number of occurrences, or the leftmost and rightmost occurrence of $P$ in $O(|P|)$ time. The locus of a string $P$ is the minimum depth node $v$ such that $P$ is a prefix of $s_v$. First we find the locus by walking downwards in the suffix tree, matching each character in $P$ in worst-case constant time using the dictionary. Once we have found $v$ we can report all the occurrences in $[\alpha, \beta]_v$ in $O(\text{occ})$ time. Alternatively, we can find the rightmost occurrence of $P$ in constant time by doing a range maximum query on the range $[\alpha, \beta]_v$ in $L$, which returns the $j \in [\alpha, \beta]$ maximizing the string position $L[j]$. We can also find the leftmost occurrence by doing a range minimum query.

Finally, note that it is possible to deamortize algorithms with expected running time using the standard technique of distributing the work evenly. Specifically, if an algorithm runs in expected $\lambda n$ time we can do $\lambda$ work for $n-1$ steps; by linearity of expectation only expected $\lambda$ work remains for the last step.

## 3 The Timely SSWSI Problem

Here we present a solution for the timely variant that matches the bounds in Theorem 1 in expectation. Section 5 shows how to get the bounds with high probability. Throughout this section we assume without loss of generality that $w$ is a power of two. Section 3.3 briefly mentions how to generalize to arbitrary $w$.

The main idea is as follows. We maintain a suffix of $S$ of length at least $w$. This suffix is segmented into at most log $w$ segments whose sizes are distinct powers of two, in increasing order from right to left. The length of the suffix we store is at most $2^0 + \ldots + 2^{\log w} = 2w - 1$. When a new character arrives, we append a new size-one segment to our data structure and merge equally-sized segments until they all have distinct sizes again. We also discard the largest segment when it no longer intersects the window. For each segment we store a suffix tree, and for every pair of adjacent segments we store a boundary tree approximately covering them both (see below). To support queries we query the suffix tree for each individual segment, and also each boundary tree. For the segments larger than the pattern, the boundary trees are sufficient to find the occurrences crossing the respective boundary. The remaining trees cover a suffix of $S$ that is $O(|P|)$ long, and we grow a suffix tree at query time to find the remaining occurrences in this suffix.
3.1 Data Structure

At any point, the data structure contains a suffix $s$ of length $w \leq |s| \leq 2w - 1$ and a segmentation of $s$ into at most $\log w$ segments. Specifically, if $|s| = 2^{b_1} + \ldots + 2^{b_k}$ for integers $b_1 < \ldots < b_k$ then we have the segmentation $s_1, \ldots, s_k$ where $|s_i| = 2^{b_i}$, and $s$ is the concatenation of the strings $s_k, s_{k-1}, \ldots, s_1$, in that order. The set $\{b_1, \ldots, b_k\}$ is unique and corresponds to the 1-bits in the binary encoding of $|s|$. Three different configurations can be seen in Figure 1.

For each segment $s_i$ we store the suffix tree $T_i$ over $s_i$, along with a range maximum query data structure over the suffix array of $s_i$. For each boundary $(s_{i+1}, s_i)$ we store the boundary tree $B_i$, which is the suffix tree over the substring centered at the boundary and extending $|s_i|$ characters in both directions. We augment $B_i$ with an additional data structure that we will use for reporting occurrences across the boundary. Let $BL_i$ be the suffix array corresponding to $B_i$. We define the modified suffix array $BL'_i$ as

$$BL'_i[j] = \begin{cases} BL_i[j] & \text{if } BL_i[j] \text{ corresponds to a suffix starting in } s_{i+1} \\ -\infty & \text{if } BL_i[j] \text{ corresponds to a suffix starting in in } s_i \end{cases}$$

We store a range maximum query data structure over $BL'_i$. Each of the data structures use $O(s_i)$ space, so the whole data structures uses $O(s) = O(w)$ space.

We note a few properties of the data structure. Let $S[n]$ be the most recent character to arrive and let $W_n = S[n - w + 1, n]$ be the current window. Then $W_n$ is a suffix of $s$ since $|s| \geq w$. The largest, and leftmost, segment $s_k$ always has size $2^{\log w} = w$; it is not larger since $\log w$ bits are sufficient to represent $|s| \leq 2w - 1$, and it is always there since $|s| \geq w$ cannot be represented with $\log w - 1$ bits. For the same reason, $s_k$ always intersects at least partially with $W_n$, and each of $s_1, \ldots, s_{k-1}$ are fully contained in $W_n$.

3.2 Queries

The idea is as follows, as exemplified in Figure 2. Any occurrence of a pattern $P$ that is fully contained in a segment is found using the suffix tree over that segment. Similarly, any occurrence that only crosses a single boundary far enough away from the end of the window is found in the respective boundary tree. Note that in the leftmost segment we must be careful to not report any occurrences that start before the left window boundary. The remaining occurrences are not contained in any of the trees in the data structure (either because they cross multiple boundaries or because they cross a single boundary $(s_{i+1}, s_i)$
We now show how to report the occurrences of P as follows, as illustrated in Figure 3. Let $s_i$ be the smallest segment where $|s_i| \geq m$. Consider any boundary $(s_{j+1}, s_j)$ to the left of $s_i$, i.e., where $j \geq i$. Since both of these segments have size at least $|s_i| \geq m$, the boundary tree $B_j$ extends at least $m$ characters in both directions from the boundary. Therefore, all the occurrences of P crossing the boundary are contained in $B_j$, and none of them can cross another boundary as well. Now
consider the suffix $R$ of $s$ containing the $m - 1$ last characters of $s_i$ and extending to the end of $s$. This substring contains all the other boundary-crossing occurrences. Furthermore, all the occurrences in $R$ cross at least one boundary since the longest consecutive part of a single segment in $R$ is the $m - 1$ characters in $s_i$. Note that the length of $R$ is at most $m - 1 + |s_{i-1}| + |s_{i-1}|/2 + \ldots + 1 < m - 1 + 2|s_{i-1}| < 3m$ since $|s_{i-1}| < m$. Thus, the number of boundary-crossing occurrences of $P$ equals the number of occurrences in $R$ plus the number of occurrences crossing the boundaries $(s_k, s_{k-1}), (s_{k-1}, s_{k-2}), \ldots, (s_{i+1}, s_i)$.

The algorithm for finding the occurrences in the sufficiently large boundary trees is as follows. Fix a boundary $(s_{k+1}, s_k)$. We match each character of $P$ in $B_x$ as it arrives. When the last character arrives we know if $|s_x| \geq m$, and also the range $[\alpha, \beta]$ corresponding to the occurrences of $P$ in the boundary tree. If $|s_x| \geq m$ (hence $x \geq i$) we report the occurrences as follows. As above we do a range maximum query to find the $j$ maximizing $BL'_x[j]$. If $BL'_x[j] = -\infty$ then all occurrences of $P$ start in $s_x$, and there are no occurrences crossing the boundary. Otherwise, $BL'_x[j]$ corresponds to the starting position of the rightmost occurrence of $P$ in $s_{x+1}$. Since all of $P$ has arrived and we now know $m$, we know that this occurrence crosses the boundary if and only if $BL'_x[j] \geq |s_x| - m + 2$ (recall that $B_x$ extends $|s_x|$ characters in both directions from the boundary). If it does not cross the boundary, then none of the other occurrences do either. Otherwise we report $BL'_x[j]$ and recurse on $[\alpha, j - 1]$ and $[j + 1, \beta]$ to find the remaining occurrences. Matching $P$ in all boundary trees takes $O(\log w)$ overall time per character, and reporting each occurrence with range maximum queries takes constant time.

We now show how to find the occurrences of $P$ in $R$ with the same bounds. Assume that we know that $2^\ell \leq m < 2^{\ell+1}$ for some integer $\ell$. We build the suffix tree over the last $3 \cdot 2^{\ell+1}$ characters of $s$, deamortized over receiving the first $2^{\ell-1}$ characters of $P$. Over the next $2^{\ell-1}$ characters we match $P$ in the tree, at a rate of two characters per new character from $P$. Then, when the $2^{\ell}$th character arrives, we have caught up to the stream $P$, and we match the remaining $m - 2^\ell$ characters as they arrive. When the last character arrives we have matched $P$ in a tree of size at least $3m$, and we can start reporting occurrences. Note that we are overestimating the size of the tree, and it potentially includes some occurrences of $P$ that are contained in $s_i$. To avoid reporting these, we also build a range maximum query data structure over the suffix array such that we can use recursive range maximum queries. When deamortized, we construct the tree in expected constant time per character of $P$. Matching $P$ also takes constant time per character. We know that $m \leq w$, so we run this algorithm simultaneously for each of the $\log w$ different choices for $\ell$, using expected $O(\log w)$ time per character in $P$. Note that the trees use $O(w)$ space in total since the sum of the space is a geometric sum where the largest term is $O(w)$.

### 3.3 Amortized Updates

We show how to support updates in amortized $O(\log w)$ time. Let $S[n]$ be the last character to arrive and as in the description of the data structure let $b_1 < b_2 < \ldots < b_k$ be the positions of the 1-indices in the binary encoding of $|s|$. When the new character $c = S[n + 1]$ arrives, we update $s$ and the segmentation $s_1 \ldots s_k$ to create the new suffix $s'$ with the new segmentation $s'_1 \ldots s'_k$. See Figure 1 for an example.

If $|s| < 2w - 1$ then we set $s' = sc$. The segmentation of $s'$ corresponds to the unique binary encoding of $|s'| = |s| + 1$, so we update the segmentation analogously to a “binary increment”. One way to do so is as follows. We create a new segment of size one over $c$. If there was not already a segment of size one, then we add the new segment and we are done. Otherwise we merge (see below) the two size-one segments to create a segment of size
two. The process cascades until we reach a size $2^b$ that does not exist in the segmentation of $s$ (i.e., the smallest index $b \not\in \{b_1, \ldots, b_k\}$). At this point we replace all of the segments $s_{i-1}, \ldots, s_1$ with $s'_1$ covering the last $2^b$ characters of $s'$. The remaining segments for $s'$ are the same as the segments $s_{b+1}, \ldots, s_k$. If $|s| = 2w - 1$ then there is a segment of each size $2^0, 2^1, \ldots, 2^\log w$. Since the segments have decreasing size from left to right, the $\log w - 1$ rightmost segments cover the last $2^0 + \ldots + 2^\log w - 1 = w - 1$ characters of $s$. Thus, after $c$ arrives, the leftmost segment of size $2^\log w = w$ no longer intersects the window. We remove it by setting $s' = s[w+1, \log w - 1]$ and update the segmentation as above.

Let $s_a$, $s_b$ and $s_c$ be three adjacent segments, in that order. To merge $s_b$ and $s_c$ we combine them into a new segment $s_d$ that spans them both, construct the suffix tree over $s_d$, and construct a range maximum query data structure on the suffix array of $s_d$. Furthermore, since $s_a$ and $s_d$ are now adjacent we also construct the boundary-spanning suffix tree for the boundary $(s_a, s_d)$ that extends $|s_d|$ characters in each direction. The construction of all of these data structures takes expected $O(|s_d|)$ time (see Section 2). Thus, it takes expected constant time per character every time it moves into a new, larger segment. Each character is contained in at most $\log w$ segments before it leaves the window, so the amortized update time is expected $O(\log w)$ per character.

Note that all but the last merge are unnecessary to actually compute $s'_1$; in the amortized setting we can simply determine where the cascade will end and immediately construct the suffix tree over the corresponding segment. However, the cascading merges will come into play in the deamortized variant.

Also note that if $w$ is not a power of two we can use a similar scheme where we allow either two simultaneous trees of size $2^{\log w}$, or one tree of size $2^\log w$. In both cases, there are some straightforward edge cases for when to remove the leftmost segment.

### 3.4 Deamortized Updates

We now show how to deamortize the updates. Unfortunately the previous construction cannot be directly deamortized since the suffix tree construction algorithm by Farach-Colton et al. [12] requires access to the whole string. Therefore, if a new character $c$ causes a cascade of merges resulting in a new segment of size $2^i$ we have to build the suffix tree over that segment when $c$ arrives.

Instead, we modify the structure slightly. When two segments of size $2^i$ become adjacent we temporarily keep both while deamortizing the cost of merging them over the next $2^i$ characters of $S$, doing expected constant work per character. Note that queries are unaffected.
The main idea is as follows. As before, we maintain suffix trees of exponentially increasing sizes, although only the $O(\log w)$ largest of them. As a result there are fewer trees to query, but also an uncovered suffix of size $\Theta(\delta)$ of the window for which we do not have any suffix trees. As in Section 3 we denote the part of $S$ covered by suffix trees by $s$ and we denote the uncovered suffix by $t$. As above, $s$ is segmented into $s_1, \ldots, s_k$.

We will first explain how to solve the problem when all patterns are long, that is, $|P| > \delta/4$, and then when all patterns are short, that is, $|P| \leq \delta/4$. Finally we show how to combine these solutions. When all the patterns are long we can afford to construct, at query time,
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a suffix tree covering \( t \). On the other hand, when all the patterns are short we can do both updates and queries in an offline fashion; we buffer queries and updates until we have approximately \( \delta/2 \) operations to do, at which point we can afford to construct a suffix tree over \( t \) in a deamortized manner. See Figure 4 for an example.

Throughout this section we assume without loss of generality that \( \delta \) is a power of two. Otherwise we instead use a more restrictive delay of \( \delta' = 2^{\lfloor \log \delta \rfloor} \) and achieve the same asymptotic bounds.

4.1 Long Patterns

We first show how to support queries if all patterns have a length \( m > \delta/4 \). We modify the data structure from Section 3 slightly. The smallest tree now has size \( \delta/2 \) as opposed to 1, so there are \( \Theta(\log w - \log(\delta/2)) = O(\log(w/\delta)) \) segments and boundary trees. The uncovered suffix \( t \) has length at most \( \delta \).

We answer queries the same way as in Section 3.2, with only small modifications. Let \( P \) be a pattern of length \( m > \delta/4 \). As before, let \( s_i \) be the smallest and rightmost segment with \( |s_i| \geq m \). We find any occurrence within a segment or crossing a single boundary by using the suffix trees over each segment and the boundary trees to the left of \( s_i \), as before. The remaining occurrences we again find by growing suffix trees of exponentially increasing sizes from the right window boundary. The only change is that we now grow the trees faster, as we must also cover \( t \), and we can afford to let the smallest tree have size \( \delta \) since we have \( m > \delta/4 \) characters in the pattern to deamortize the work over. As above, let \( R \) be the string covering the \( m-1 \) last characters of \( s_i \) and extending to the right window boundary, which now also includes \( t \). As \( |t| < \delta \) the length of \( R \) is \( |R| < 3m + \delta < 7m \). Assuming \( 2^\ell \leq m < 2^{\ell+1} \), we build the suffix tree of size \( 7 \cdot 2^{\ell+1} \) and match \( P \) in it, amortized over the characters of \( P \). As we have \( m > \delta/4 \) characters to deamortize the work over, we only do this for each choice of \( \ell \) where \( 2^{\ell+1} \geq \delta \), which results in \( O(\log w - \log \delta) = O(\log(w/\delta)) \) work per character in \( P \). As in Section 3.2 we use recursive range maximum queries to avoid double reporting any occurrences of \( P \) that are also in \( s \). As there are also only \( O(\log(w/\delta)) \) segments and boundary trees we spend \( O(\log(w/\delta)) \) time per character in \( P \). Note that we answer these queries without delay.

Updates are performed as follows. For each segment of \( \delta/2 \) characters that arrives we construct the suffix tree over it, deamortized over the next \( \delta/2 \) characters of \( S \). We merge suffix trees as before, also deamortized over new characters of \( S \). The induction proof from Section 3.4 still works by modifying the base case; the merging of two trees of size \( \delta/2 \) takes \( \delta/2 \) characters, at which point another tree of size \( \delta/2 \) is constructed. The inductive step follows from the fact that \( \delta \) is a power of two. Thus, we spend expected \( O(\log(w/\delta)) \) time per update.

4.2 Short Patterns

We now show how to support queries if all patterns have a length \( m \leq \delta/4 \). We extend the data structure with a buffer of size \( \delta \). This buffer will contain queries that we have not yet answered and characters for \( S \) that we have not yet processed. The total space is still \( O(w + \delta) = O(w) \).

Whenever a character from \( S \) arrives we append it to both \( t \) and to the buffer. When a pattern arrives we append the full pattern to the buffer, and along with it we store the current position of the right window boundary. Once the buffer has more than \( \delta/2 \) characters
We immediately allocate a new buffer of size $\delta$ and flush the old buffer as follows. Note that at this point there are strictly less $\frac{3}{4}\delta$ characters in the buffer since each pattern is short.

When we flush the buffer, we first answer all the buffered queries, and then we process all the buffered updates. We deamortize this work over the next $4\log(w/\delta)$ characters that arrive from either stream. To answer the buffered queries we do as follows. Let $P_1, \ldots, P_t$ be the patterns in the buffer, let $m_i = |P_i|$, and let $M = \sum_{1 \leq i \leq t} m_i$. We have $M < \delta$. We start by building a suffix tree over $t$, along with a range maximum query data structure over the suffix array of $t$. This takes expected $O(\delta)$ time. An occurrence of $P_i$ is either contained in $s$, or it crosses the boundary $(s, t)$, or it is contained in $t$. Since $P_i$ is smaller than each segment $s_j$ we can find all the occurrences within $s$ using the suffix trees over the segments and the boundary trees in $O(m_i \log(w/\delta))$ time. To find the occurrences crossing the boundary we build the KMP matching automaton [20] for $P_i$. In it we match the string that is centered at the boundary $(s, t)$ and extends $m_i - 1$ characters in each direction. This takes $O(m_i)$ time. To find the occurrences in $t$ we match $P_i$ in the suffix tree over $t$ in $O(m_i)$ time. In total, this takes $O(M \log(w/\delta)) = O(\delta \log(w/\delta))$ time for all the patterns, or expected $O(\log(w/\delta))$ time per character when deamortized. Note however, that after $P_i$ arrived more characters from $S$ could have arrived and been appended to $t$. We must therefore take care not to report any occurrences of $P_i$ that extend past what was the right window boundary when $P_i$ arrived. The KMP automaton finds the occurrences in left-to-right order, and in $t$ we avoid reporting too far right using recursive range minimum queries.

Finally, we process each update in the buffer in the order they arrived, using the same procedure as for long patterns. This takes $O(\log(w/\delta))$ time per update and $O(\delta \log(w/\delta))$ time in total. Thus flushing the buffer takes expected $O(\log(w/\delta))$ time per character since we deamortize the expected $O(\delta \log(w/\delta))$ work over $\delta/4$ characters. Since we allocate a new buffer immediately when we begin flushing, we will complete the flush before the next flush begins.

4.3 Both Long and Short Patterns

We now show how to combine the solutions for short and long patterns, to obtain a solution that handles patterns of any length. The data structure is the same as for small patterns above. As above, we append each new character to the buffer. However, whenever we start
streaming a pattern we also proceed as if $P$ were long. If $P$ turns out to fit in the buffer without triggering a flush (which might also happen if $P$ is long), we simply discard the work we did for the long-pattern case. However, if adding $P$ to the buffer results in more than $\frac{d}{2}$ characters being in the buffer, then $P$ must be long. We immediately start flushing the buffer (ignoring the characters related to $P$) and also continue processing $P$ as a long pattern. Note that since we are potentially streaming a long pattern while batch processing the updates in the buffer, the data structure might change while we are matching in it. However, it only changes when a merge finishes, replacing a pair of suffix trees by a larger tree. If this happens we keep the old trees in memory until we are done processing the pattern, at which point we discard them.

We obtain the following theorem.

Theorem 4. Let $S$ be a stream and let $w \geq 1$ and $\delta \geq 1$ be integers. We can solve the $(w, \delta)$-SSWSI problem on $S$ with an $O(w)$ space data structure that supports Update and Report in expected $O(\log(w/\delta))$ time per character. Furthermore, Report uses additional worst-case constant time per reported occurrence.

5 Obtaining High Probability

In this section we show how to improve the time bounds to $O(\log(w/\delta))$ with probability $1 - w^{-d}$ for any constant $d \geq 1$.

The expectation in the time bounds in Section 4 comes from the construction of suffix trees (recall that we also build suffix trees at query time). Below, in Lemma 5, we prove that given a string $K$ of length $k = O(w)$ we can construct the suffix tree over $K$ in $O(k)$ time with probability $1 - 1/w^{1+\epsilon}$, using additional $O(w/\log w)$ space. We use this algorithm to construct suffix trees during updates and queries, deamortizing them as before and doing $O(\log(w/\delta))$ work per character that arrives. When a new character arrives from $S$ or $P$, at most $O(\log(w/\delta)) = O(\log w)$ suffix tree constructions will finish. At this point, we finish constructing those trees that did not finish in time, that is, used more more time than what was allotted to them. By the union bound, the probability that any of them fail to finish in time (and thus incurring extra construction cost) is no more than $c \log w/w^{1+\epsilon}$ for some constant $c$ which is no more than $1/w$ for large $w$. Thus, for each character from $S$ or $P$ we spend $O(\log(w/\delta))$ time with high probability in $w$. We obtain the $1 - 1/w^{d}$ probability bound by probability boosting, running $d = O(1)$ independent copies of the construction algorithm simultaneously. The algorithm from Lemma 5 uses additional $O(w/\log w)$ space, but we are never constructing more than $O(\log w)$ suffix trees, so the space usage is $O(w)$ in total.

Furthermore, as mentioned in Section 2, we previously used an FKS dictionary [15] to store the edges to support reporting queries in worst-case constant time per character in the pattern. The construction time of this dictionary is expected linear, so it can no longer be used. Instead we use a dictionary by Dietzfelbinger and Meyer auf der Heide [11]. If there are $n$ elements in the dictionary it supports searches in worst-case constant time and any sequence of $\frac{1}{2}n$ updates takes constant time per update with probability $1 - 1/w^{d}$ for any constant $d \geq 1$. We store all the edges of all the suffix trees in one such dictionary. At all times, we keep $\Theta(w)$ dummy-elements in the dictionary to ensure that we get good probability bounds in terms of $w$, and we choose $d'$ large enough that any sequence of $O(w)$ operations (e.g., the construction of any one of our suffix trees) runs in $O(w)$ time with probability $1 - 1/w^{d+\epsilon}$.
Universal Hashing

Before we prove Lemma 5 we restate some basic facts about universal hashing, introduced by Carter and Wegman [9]. Let $M, m > 0$ be integers, $\mathcal{H}$ be a set of functions $[0,M] \to [0,m]$, and $h \in \mathcal{H}$ be selected uniformly at random. Then $\mathcal{H}$ is universal if $P[h(x) = h(y) \mid x \neq y] \leq 1/m$. Let $R \subseteq [0,M]$ and $|R| = r$. It follows from the union bound that $h$ has a collision on $R$ with probability at most

$$P[h(x) = h(y) \text{ for some } x \neq y] \leq \sum_{x \neq y \in R} P[h(x) = h(y)] = \frac{r(r-1)}{2} \cdot \frac{1}{m} < \frac{r^2}{m}. \quad (1)$$

In particular, if $m = r^c$ for constant $c \geq 1$ then $h$ is injective (i.e., has no collisions) on $R$ with probability at least $1 - 1/r^{c-2}$. Carter and Wegman gave several classes of universal hash functions from which we can sample a function uniformly at random in constant time.

Fast Suffix Tree Construction

We now prove Lemma 5, showing how to construct our suffix trees in linear time with high probability.

**Lemma 5.** Given a string $K$ of length $k \leq 2w$ there is an algorithm that uses $O(k+w/\log w)$ space and constructs the suffix tree over $K$ in $O(k)$ time with probability $1 - 1/w^{1+\epsilon}$ for some $\epsilon > 0$.

**Proof.** Let $\sigma = \{K[i] \mid i \in [1,k]\} \subseteq \Sigma$ be the alphabet of $K$. We show how to, in $O(k)$ time, find a function $h : \Sigma \to [1,k^{O(1)}]$ such that $h$ is injective on $\sigma$ with probability at least $1 - 1/w^{1+\epsilon}$. If $h$ is injective on $\sigma$, we can construct the suffix tree over $K'$ where $K'[i] = h(K[i])$ in time $O(sort(k,k^{O(1)})) = O(k)$ using radix sort. After the tree is constructed we can substitute for the original alphabet in linear time. Therefore, the construction algorithm finishes in $O(k)$ time with probability at least $1 - 1/w^{1+\epsilon}$ (otherwise we make no guarantee on the construction time and we can build the suffix tree in any way).

For some $m$ to be determined later, let $f : \Sigma \to [1,m]$ be chosen uniformly at random from a class of universal hash functions. By Equation 1, the probability that $f$ has a collision on $\sigma$ is

$$P[f \text{ has collisions on } \sigma] < \frac{|\sigma|^2}{m} < \frac{k^2}{m}.$$  

We divide into the cases of large trees ($k \geq w^{1/5}$) and small trees ($k < w^{1/5}$). If $k$ is large then $w^{1/5} \leq k \leq 2w$, and we set $m = w^4$ so the probability that $f$ has a collision is at most

$$\frac{k^2}{m} \leq \frac{(2w)^2}{w^4} = \frac{4}{w^2} \leq \frac{1}{w^{1+\epsilon}}$$

for some $\epsilon > 0$. We check whether $f$ is injective by sorting the set $\{(x,f(x)) \mid x \in \sigma\}$ with respect to the $f(\cdot)$-values and checking if two consecutive elements $(x_1,f(x_1))$ and $(y,f(y))$ have $x_1 \neq y$ and $f(x_1) = f(y)$. This takes time $O(sort(k,w^4)) = O(k)$ using radix sort since $k \geq w^{1/5}$. If $f$ is injective we set $h = f$, concluding the proof of the large case.

If $k$ is small then we allocate an array $A$ of length $w/\log w$ in constant time. For simplicity we assume that $A$ is initialized such that $A[i] = 0$ for all $i$. This can be avoided using standard constant-time initialization schemes; assume each entry in $A$ contains an arbitrary value initially. We maintain two other arrays $B$ and $C$ such that if we have written a value to $A[i]$
at least once then $A[i]$ is a pointer to some $B[j]$, $B[j]$ is a pointer to $A[i]$, and $C[j]$ stores the value most recently written to $A[i]$. From this we can determine if $A[i]$ has been initialized (check if the pointers match), and if it has not we can initialize it in constant time.

Then we set $m = w/\log w$ such that the probability that $f$ has a collision is no more than

$$\frac{k^2}{m} < \frac{w^{2/5}}{w/\log w} = \frac{\log w}{w^{3/5}} = \frac{\log w}{w^{1/2}} \cdot \frac{1}{w^{1/10}} \leq \frac{1}{w^{1/10}}$$

for $w \geq 16$. We check if $f$ is injective on $\sigma$ by for each character $x$ in $K$ setting $A[f(x)] = x$ and seeing if two distinct characters hash to the same index. If $f$ is injective we then arbitrarily assign the values $1, \ldots, |\sigma|$ to the now non-zero indices of $A$ and let $h(x) = A[f(x)]$ (at this point we know $\sigma$ since it is equal to the number of entries in $A$ that we modified). To boost the probability of success we run this algorithm up to eleven times with independent choices for $f$. The probability that all of them fail is at most $1/w^{11/10} \leq 1/w^{1+\epsilon}$ concluding the proof for the small case.

In conjunction with Theorems 3 and 4, this proves Theorem 1.

6 Conclusion and Future Work

We have studied two variants of the streaming sliding window string indexing problem; the timely variant, where queries must be answered immediately, and the delayed variant where a query may be answered at any point within the next $\delta$ characters received, for a specified parameter $\delta$. For a sliding window of size $w$ we have given an $O(w)$ space data structure that, in the timely variant, supports updates in $O(\log w)$ time with high probability and queries in $O(\log w)$ time with high probability per character in the pattern; each occurrence is reported in additional constant time. For the delayed variant we improved these bounds to $O(\log(w/\delta))$, where each occurrence is still reported in constant time.

One open problem is whether these bounds can be improved. Another is to find efficient solutions when queries may be interleaved with new updates to the stream. That is, while you are streaming a pattern, new characters of $S$ might arrive that move the current window.

References


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Faster Algorithms for Computing the Hairpin Completion Distance and Minimum Ancestor

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Abstract

Hairpin completion is an operation on formal languages that has been inspired by hairpin formation in DNA biochemistry and has many applications especially in DNA computing. Consider $s$ to be a string over the alphabet \{A, C, G, T\} such that a prefix/suffix of it matches the reversed complement of a substring of $s$. Then, in a hairpin completion operation the reversed complement of this prefix/suffix is added to the start/end of $s$ forming a new string.

In this paper we study two problems related to the hairpin completion. The first problem asks the minimum number of hairpin operations necessary to transform one string into another, number that is called the hairpin completion distance. For this problem we show an algorithm of running time $O(n^2)$, where $n$ is the maximum length of the two strings. Our algorithm improves on the algorithm of Manea (TCS 2010), that has running time $O(n^2 \log n)$.

In the minimum distance common hairpin completion ancestor problem we want to find, for two input strings $x$ and $y$, a string $w$ that minimizes the sum of the hairpin completion distances to $x$ and $y$. Similarly, we present an algorithm with running time $O(n^2)$ that improves by a $O(\log n)$ factor the algorithm of Manea (TCS 2010).

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1 Introduction

1.1 Motivation and informal problem definition

Hairpin completion is an operation on formal languages that has been inspired by hairpin formation in DNA biochemistry and has many applications especially in DNA computing [11, 12, 14, 15]. This operation has been inspired by three biological principles: Watson-Crick complementarity, DNA annealing and DNA lengthening through polymerases. The DNA chain is a molecule consisting of two intertwined strands, each strand being composed by nucleotides: A(Adenine), C(cytosine), G(guanine) and T(thymine). The two strands which
form the DNA molecule are kept together by the hydrogen bond between the bases: A bonds with T and C with G. This paradigm is usually referred to as the Watson-Crick complementarity [25].

Another important bio-chemical principle is annealing, the process of fusing two single stranded molecules by complementary base. DNA lengthening through polymerases is a phenomenon that produces a complete double stranded DNA molecule as follows: one starts with two single strands such that one (called primer) is bonded to a part of the other (called template) through Watson-Crick complementarity and a polymerization buffer with many copies of the four nucleotides. The polymerases will then concatenate to the primer by complementing the template [22].

We now begin to informally explain the hairpin completion operation and how it can be related to the biological concepts presented above. Consider $s$ to be a string over the alphabet \{A,C,G,T\} such that a prefix/suffix of it matches to the reversed complement of a substring of $s$. Then, the reversed complement of this prefix/suffix is added to the beginning/ending of $s$ forming a new string as can be visualized in Figure 1. The mathematical expression of this hypothetical situation defines the hairpin completion operation. Starting with a single string, one can generate a set of strings using this formal operation: via hairpin completion, a new string can be created for each possible pairing between a prefix or suffix and a complementary substring. In addition, one could be interested in knowing how many iterations of hairpin completion are required to transform one string into another. In this way, the hairpin completion distance between two strings was defined as the minimum number of times we must iterate the hairpin completion operation, starting from one of the two string, in order to obtain the other. Further, one can also be interested in finding for two strings, a common ancestor that minimizes the sum of the hairpin completion distances to those strings. This ancestor is called minimum distance common hairpin completion ancestor.

![Figure 1](image-url) An illustration of the left and right hairpin completion operations.

### 1.2 Previous and related work

The hairpin completion operation has been introduced by Cheptea, Martin-Vide and Mitrana [4]. In several papers, the hairpin completion and other familiar operations have been studied [3, 5, 6, 7, 8, 9, 13, 17, 19, 20, 21, 22, 23, 24].

Hairpin reduction [3, 22, 23] was introduced as an inverse operation for hairpin completion. The hairpin reduction of a string $x$ consists of all strings $y$ such that $x$ can be obtained from $y$ by hairpin completion. Further, two variants of hairpin completion were considered, as they seem more appropriate for practical implementation: hairpin lengthening and bounded hairpin completion [9, 19, 21]. The first variant consist of adding a prefix or a suffix of $\gamma$. The second variant assumes that the length of the added prefix or suffix is bounded by a constant. Besides the algorithmic aspects, hairpin completion operation has been studied from the language theory point of view in several papers [5, 6, 8, 13, 17].

Manea and Mitrana introduced the minimum distance common $k$-hairpin completion ancestor of two strings in [22] where they presented a cubic time algorithm to compute the ancestor. Afterwards, Manea, Martin-Vide, and Mitrana [20] suggested a cubic time
algorithm to tackle the $k$-hairpin completion distance problem. In addition, in [18] improved the time complexity to $O(n^2 \log n)$ to both problems, where $n$ is the length of the longest string.

1.3 Our results

The focus of this paper is on two algorithmic problems related to iterated hairpin completion: $k$-hairpin completion distance and minimum distance common $k$-hairpin completion ancestor. Our main results are improving the upper bound on both problems with a $\log n$ factor, from $O(n^2 \log n)$ to $O(n^2)$. For the $k$-hairpin completion distance, our speedup is based on using incremental tree, a data structure proposed by Kaplan and Shafrir [10] which can support in constant time the following operations in a weighted tree: return the edge with minimum weight on a path and add a leaf to the tree. Our algorithm for finding a minimum distance $k$-hairpin completion ancestor of two strings $(x, y)$ is based on dynamic programming technique presented in [18]. As in [18], we are interested in constructing the table $DP_x$, where $DP_x[i][j]$ represents the minimum number of $k$-hairpin completion operations to transform $x[i \ldots j]$ into $x$. Similarly, we would like to compute a table $DP_y$. Our speedup relies in an $O(n^2)$ time algorithm for computing these tables by rephrasing the problem of computing $DP_x$ in terms of shortest distances in a graph and replacing the segment tree used in [18] with doubly linked list and changing the order we process the cells in the matrix.

2 Preliminaries

We start with basic notations related to strings. An alphabet $\Sigma$ is a finite, non-empty set of symbols. Throughout this paper, we mostly discuss strings over the alphabet $\Sigma = \{A, C, G, T\}$. For a letter $x \in \Sigma$, we denote as $\pi$ the letter in $\Sigma$ that is complementary to $x$. For the previously mentioned alphabet, we have $\pi = T$ and $\pi = G$. The set of all strings over an alphabet $\Sigma$ is denoted by $\Sigma^*$. The empty string is denoted as $\lambda$, and $\Sigma^* = \Sigma^* \setminus \{\lambda\}$. Given a string $w \in \Sigma^*$, we denote by $|w|$ its length. If $w = xy$, $x, y \in \Sigma^*$ then $x$ is called prefix and $y$ a suffix. For a string $w$, $w[i \ldots j]$ denotes the substring of $w$ starting at position $i$ and ending at position $j$, $1 \leq i \leq j \leq |w|$. Given a string $s \in \Sigma^*$, we denote by $\overline{s} = s_1 \overline{s_2} \ldots \overline{s_{|s|}}$ the complement of the string $s$ and $s^{R}$ the reversed string of $s$, i.e. $s^R = s_{|s|} s_{|s|-1} \ldots s_1$.

Incremental tree is a data structure introduced by Kaplan and Shafrir [10] based on a similar structure of Alstrup and Holm [1] for the level ancestor problem, to maintain a rooted tree $T$, with an integer weight on each edge, such that the following operations are supported in $O(1)$ amortized time:

- add-leaf$_T(v, w, c)$: Add a new leaf $v$ with parent $w$ to $T$. The weight of the edge $(v, w)$ is $c$.
- add-root$_T(v, c)$: Add a new root $v$ to $T$. The old root $(r)$ becomes a child of $v$ and the weight of edge $(r, v)$ is $c$.
- min-edge$_T(v, w)$: Returns the edge with minimum weight on the path from $v$ to $w$.
- change-weight$_T(v, c)$: $v$ is a leaf or $v$’s parent is the root of $T$. Changes the weight of the edge between $v$ and its parent to $c$.

From this data structure we will use just add-leaf$_T$ and min-edge$_T$ operations.

2.1 Hairpin Operations

For a string $x \in \Sigma^+$ and a positive integer $k \in \mathbb{N}$, $k$-hairpin completion is a family of transformations that can be applied to $x$. When applying a left $k$-hairpin completion, we select a non-empty suffix $\gamma$ of $x$ such that $x$ can be partitioned into $x = \alpha \beta \alpha^R \gamma$ with $\alpha, \beta \neq \lambda$ and $\beta \neq \lambda$. The process involves adding a hairpin to $x$, resulting in $x = \alpha \beta \alpha^R \gamma$.

Furthermore, the focus of this paper is on two algorithmic problems related to iterated hairpin completion: $k$-hairpin completion distance and minimum distance common $k$-hairpin completion ancestor.
\(\alpha, \beta, \gamma \in \Sigma^+\) and \(|\alpha| = k\). We execute the left hairpin operation by appending \(\gamma \overline{\alpha} R\) to the beginning of \(s\). Formally, the set of strings that can be obtained from \(s\) by applying a single left \(k\)-hairpin operation is denoted as

\[
HCL_k(x) = \{\gamma \overline{\alpha} R | x = \alpha \beta \gamma, |\alpha| = k, \alpha, \beta, \gamma \in \Sigma^+\}
\]

A right \(k\)-hairpin completion is defined in a symmetrical manner and the set of strings that can be obtained from \(s\) by applying a single right \(k\)-hairpin completion operation is denoted as

\[
HCR_k(x) = \{x \gamma \overline{\alpha} R | x = \gamma \alpha \beta \gamma, |\alpha| = k, \alpha, \beta, \gamma \in \Sigma^+\}
\]

**Example 1.** The string \(s = GAATCT\) can be partitioned into \(\alpha = GA, \beta = A, \overline{\alpha} R = TC\) and \(\gamma = T\). Applying the left hairpin completion operation on \(s\) with this partitioning yields the string \(AGAATCT\). Also, \(s\) can be partitioned into \(\gamma = GA, \alpha = A, \beta = TC, \overline{\alpha} R = T\) and by applying right hairpin completion operation we obtain \(GAATCTTC\).

Collectively, the set of strings that can be obtained from \(x\) either by applying a right or a left \(k\)-hairpin completion operation is denoted as

\[
HC_k(x) = HCL_k(x) \cup HCR_k(x)
\]

The hairpin completion is the variant of the \(k\)-hairpin completion where we do not place a bound on the length of prefix. The hairpin completion of \(x\) is defined by:

\[
HC(x) = \bigcup_{k \geq 1} HC_k(x)
\]

We extend the notation of hairpin completion to sets of strings in the following way, for a set \(L \subseteq \Sigma^*\) and a positive integer \(k\),

\[
HC_k(L) = \bigcup_{x \in L} HC_k(x) \quad HC(L) = \bigcup_{x \in L} HC(x)
\]

For every non negative integers \(k, i\) and string \(x \in \Sigma^+\), we denote as \(HC^i_k(x)\) the set of strings that can be obtained from \(x\) using exactly \(i\) \(k\)-hairpin completion operations and \(HCA^i_k(x)\) as the set of strings that are obtainable from \(x\) using any number of \(k\)-hairpin completion operations. Similarly, we denote as \(HC^i(x)\) and \(HCA^i(x)\) the sets of strings obtainable from \(x\) by applying \(i\) (resp. any number) of hairpin operations, respectively. Formally,

\[
HC^0_k(x) = \{x\} \quad HC^{i+1}_k(x) = HC_k(HC^i_k(x)) \quad HC^*_k(x) = \bigcup_{i \geq 0} HC^i_k(x)
\]

\[
HC^0(x) = \{x\} \quad HC^{i+1}(x) = HC(HC^i(x)) \quad HC^*(x) = \bigcup_{i \geq 0} HC^i(x)
\]

\[
HC^*_k(L) = \bigcup_{x \in L} HC^*_k(x) \quad HC^*(L) = \bigcup_{x \in L} HC^*(x)
\]

**Definition 2 (\(k\)-Hairpin Completion Common Ancestor).** A string \(w\) is a common \(k\)-hairpin completion ancestor of two strings \(x\) and \(y\) if \(\{x, y\} \subseteq HC^*_k(w)\). We denote the set of common \(k\)-hairpin ancestors of \(x\) and \(y\) as \(HCA_k(x, y)\).
Definition 3 (k-Hairpin Completion Distance). Given two strings \( x \) and \( y \) such that \( |x| \leq |y| \), the \( k \)-hairpin completion distance between \( x \) and \( y \) is the minimal number of \( k \)-hairpin operations required to obtain \( y \) from \( x \). Formally

\[
HCD_k(x, y) = \left\{ \begin{array}{ll}
\min\{t | x \in HCA^t_k(y)\} & \text{if } x \notin HCA^t_k(y) \\
\infty & \text{otherwise}
\end{array} \right.
\]

Definition 4 (Minimum Distance k-hairpin Completion Ancestor). For two strings \( x, y \in \Sigma^* \), a \( k \)-hairpin completion ancestor \( w \in HCA_k(x, y) \) is a minimum distance \( k \)-hairpin completion ancestor of \( x \) and \( y \) if \( \forall w' \in HCA_k(x, y) \) it holds that \( HCD_k(w, x) + HCD_k(w, y) \leq HCD_k(w', x) + HCD_k(w', y) \), i.e. \( w \) minimizes the sum of the \( k \)-hairpin completion distances from \( x \) and from \( y \).

Definition 5 (Border). Given a string \( s[1 \ldots n] \in \Sigma^+ \), Border\((s)\) is the length of the longest prefix of the string \( s \) which is also a complemented reversed suffix of this string. Formally, Border\((s)\) = \( \max\{t | s[1 \ldots 1 + t - 1] = \overline{s}[n - t + 1 \ldots n]\} \cup \{0\} \). This definition can be easily extended for any substring \( s[i \ldots j] \) in the following way: Border\((s[i \ldots j])\) = \( \max\{t | s[i \ldots i + t - 1] = \overline{s}[j - t + 1 \ldots j]\} \cup \{0\} \).

Remark 6. Note that the above definition for border is different than the common definition, which is usually the largest prefix of \( x \) which is also a suffix of \( x \).

Since in the \( k \)-hairpin completion operation we have to make sure that \( |\alpha| = k \), we introduce the definition of \( k \)-Border.

Definition 7 (k-Border). Given a string \( s \in \Sigma^+ \), k-Border\((s)\) = \( \max\{\text{Border}(s) - k, 0\} \).

Hairpin reduction is the inverse operation of hairpin completion. The hairpin reduction of a string \( x \) consists of all strings \( y \) such that \( x \) can be obtained from \( y \) by hairpin completion. For a string \( x \in \Sigma^* \) and a positive integer \( k \in \mathbb{N} \), \( k \)-hairpin reduction is a family of transformations that can be applied to \( x \). When applying a left hairpin reduction, we select a non-empty prefix \( \gamma \) of \( x \) such that \( x \) can be partitioned into \( \gamma\alpha\beta\overline{\gamma}\overline{R} \) with \( \alpha, \beta, \gamma \in \Sigma^+ \) and \( |\alpha| = k \). We execute the left hairpin reduction operation by deleting \( \gamma \). Formally, the set of strings that can be obtained from \( x \) by applying a single left \( k \)-hairpin reduction operation is denoted as

\[
HRL_k(x) = \{\alpha\beta\overline{\gamma}\overline{R} | x = \gamma\alpha\beta\overline{\gamma}\overline{R}, |\alpha| = k, \alpha, \beta, \gamma \in \Sigma^+\}
\]

A right \( k \)-hairpin reduction operation is defined in a symmetrical manner and the set of strings that can be obtained from \( x \) by applying a single right \( k \)-hairpin reduction operation is denoted as

\[
HRR_k(x) = \{\gamma\alpha\beta\overline{R} | x = \gamma\alpha\beta\overline{R}\overline{\gamma}, |\alpha| = k, \alpha, \beta, \gamma \in \Sigma^+\}
\]

The set of strings that can be obtained from \( x \) either by applying a left or a right \( k \)-hairpin reduction operation is denoted as

\[
HR_k(x) = HRL_k(x) \cup HRR_k(x)
\]

The hairpin reduction is the variant of the \( k \)-hairpin reduction where we do not place a bound on the length of prefix. The hairpin reduction of \( x \) is defined by:

\[
\text{HR}(x) = \bigcup_{k \geq 1} HR_k(x)
\]

We make the following observation.
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Observation 8. Let $x[1\ldots n]$ be a string with $k$-Border $l$.

$$HRL_k(x) = \bigcup_{j\in[1\ldots l]} \{x[j+1\ldots n]\} \quad \text{and} \quad \text{HRR}_k(x) = \bigcup_{j\in[1\ldots l]} \{x[1\ldots n-j]\}$$

$$HR_k(x) = \bigcup_{j\in[1\ldots l]} \{x[j+1\ldots n], x[1\ldots n-j]\}$$

Now we are ready to introduce the problems that we study in this paper.

Problem 1 (Hairpin completion distance). Let $\Sigma$ be the alphabet and $x, y \in \Sigma^+$. Compute $HCD_k(x, y)$.

Problem 2 (Minimum distance common hairpin completion ancestor). Let $\Sigma$ be the alphabet and $x, y \in \Sigma^+$. Compute a minimum-distance common $k$-hairpin completion ancestor of $x, y$.

2.2 Suffix Tree and Extension queries

The suffix tree [26] is a useful string data structure.

Definition 9. Let $S_1, \ldots, S_k$ be strings over alphabet $\Sigma$ and let $\$ \notin \Sigma$.

A trie of strings $S_1, \ldots, S_k$ is an edge-labeled tree with $k$ leaves. Every path from the root to a leaf corresponds to a string $S_i$ with a $\$ symbol appended to its end. The edges on this path are labeled by the symbols of $S_i$. Strings with a common prefix start at the root and follow the same path of the prefix, the paths split where the strings differ.

A compacted trie is a trie with every chain of edges connected by degree-2 nodes contracted to a single edge whose label is the concatenation of the symbols on the edges of the chain.

Let $S = S[1], \ldots, S[n]$ be a string over alphabet $\Sigma$. Let $\{S_1, \ldots, S_n\}$ be the set of suffixes of $S$, where $S_i = S[i\ldots n]$, $i = 1, \ldots, n$. A suffix tree of $S$ is the compacted trie of the suffixes $S_1, \ldots, S_n$.

For every node $u$, we call the concatenation of the labels on the path from the root to $u$ the locus of $u$ denoted as $L(u)$. For an edge $e$ in the compact trie, we use the same notation $L(e)$ to denote the label (or the locus) of $e$. Finally, for a downwards path $P$ in the compact trie, the locus $L(P)$ is the concatenation of the loci of the edges in $P$. In a compact trie, an edge $e$ can have label s.t. $|L(e)| > 1$. We refer to the symbol $L(e)[1]$ as the symbol of $e$.

Theorem 10 (Weiner [26]). For finite alphabet $\Sigma$, the suffix tree of a length-$n$ string can be constructed in time $O(n)$. For general alphabets it can be constructed in time $O(n \log \sigma)$, where $\sigma = \min(|\Sigma|, n)$.

For two strings $S[1\ldots n]$ and $T[1\ldots m]$, a string $P[1\ldots p]$ is a common prefix of $S$ and $T$ if $S[1\ldots p] = T[1\ldots p] = P$. We say that $P$ is the longest common prefix (LCP) of $S$ and $T$ if $P$ is a common prefix and $m = p$ or $n = p$ or $S[p+1] \neq T[p+1]$. Similarly, a string $A[1\ldots a]$ is a common suffix of $S$ and $T$ if $S[n-a+1\ldots n] = T[m-a+1\ldots m] = A$. $A$ is the longest common suffix (LCS) if $n = a$ or $m = a$ or $S[n-a] \neq T[m-a]$. Collectively, we refer to LCP and LCS as longest common extensions (LCE).

By preprocessing the suffix tree of a string $S$ for level ancestor queries [2], we can obtain the following.

Lemma 11 (Longest Common Extension Data Structure). A string $S$ can be preprocessed in $O(n)$ time to support the following queries in $O(1)$ time.
1. $LCP(i, j)$ - return the length of the longest common prefix of $S[i \ldots n]$ and $S[j \ldots n]$
2. $LCS(i, j)$ - return the length of the longest common suffix of $S[1 \ldots i]$ and $S[1 \ldots j]$
By constructing the above data structure for the string $x^i y x^j$ with $\not\in \Sigma$, we obtain the following.

**Corollary 12.** We can process a string $S[1 \ldots n]$ in linear time to construct a data structure for answering the following query in $O(1)$ time.

\[
\text{k-Border}(S[i \ldots j]) = \text{Return the length of the k-Border of } S[i \ldots j].
\]

## 3 Hairpin completion distance

In this section we study Problem 1.

Our algorithm is based on the dynamic programming technique presented in [18]. For the sake of clarity, we briefly describe this technique. Without loss of generality, we assume that $|x| \leq |y|$ and $n = |y|, m = |x|$. We are interested in computing a dynamic programming table $DP[i][j]$ with dimensions $n \times n$. For every two indices $1 \leq i \leq j \leq n$, we define $DP[i][j]$ to be the minimum number of k-hairpin completion operations to transform $x$ into $y[i \ldots j]$. Formally, $DP[i][j] = \text{HCD}_k(x, y[i \ldots j])$.

**Definition 13.** Given two non-negative integers $i, j$ ($1 \leq i \leq j \leq n$), $L_j$ represents the $DP$ values of all strings that can generate the substring $y[i \ldots j]$ through a single left $k$-hairpin completion operation (elements of the set $HRL_k(y[i \ldots j])$).

**Claim 14.** $L_j = \{DP[i + 1][j], \ldots, DP[i + l][j]\}$ where $l$ is the k-Border of $y[i \ldots j]$.

**Definition 15.** Given two non-negative integers $i, j$ ($1 \leq i \leq j \leq n$), $R_i$ represents the $DP$ values of all strings that can generate the substring $y[i \ldots j]$ through a single right $k$-hairpin completion operation (elements of the set $HRR_k(y[i \ldots j])$).

**Claim 16.** $R_i = \{DP[i][j - l], \ldots, DP[i][j - 1]\}$ where $l$ is the k-Border of $y[i \ldots j]$.

Correctness of Claim 14 and Claim 16 is based on Observation 8.

**Lemma 17.** Given two non-negative integers $i, j$ ($1 \leq i \leq j \leq n$), we have that $DP[i][j] = \min(\min(L_j), \min(R_i)) + 1$.

For the proof of Lemma 17 we refer to [18].

All positions in $DP$ are initialized with $\infty$. We start by considering the base cases. These are represented by all subsequences $y[i \ldots j] = x$. To determine them, we use any pattern matching algorithm which runs in linear time, for example KMP [16] and set $DP[i][j] = 0$. Analyzing the elements of the sets $L_j$ and $R_i$, it can be seen that they actually represent continuous blocks from line $i$ or column $j$. Thus, determining the minimum values from each of those sets is a range minimum query.

**Definition 18.** Given two non-negative integers $i, j$ ($1 \leq i \leq j \leq n$), $DSL_j$ represents the data structure that keeps the $DP$ values of column $j$ and $DSR_i$ represents the data structure that keeps the $DP$ values of row $i$. (Note that we don’t have to keep the values below the main diagonal)

Naively, $DSL_j$ and $DSR_i$ could be arrays, which leads to constant update time, but linear query time. The overall complexity of the algorithm with this naive implementation is $O(n^3)$. In [18], the algorithm is implemented using segment trees, which leads to a logarithmic time for queries and for updates.
We compute the $DP$ matrix in increasing order of difference $j - i$ (parallel with the main diagonal). Red line represents $DSL_i$ and the green line $DSR_i$.

Considering that the update operations are append-like, i.e. they are only done after the first/last index of $DSL_j$ and $DSR_i$, we propose using an incremental tree. The advantages of this approach consist in the fact that this structure can perform query and update operations in constant time. Practically, we keep an incremental tree for each row and column. A row or a column in the matrix represents a particular case of a tree, more precisely a chain. For the range minimum query needed in the computation of $DP[i][j]$ we use incremental tree's $\text{min-edge}_T$ operation. After we compute the $DP[i][j]$ value, we have to add to $DSR_i$ and $DSL_j$. This can be done by using the $\text{add-leaf}_T$ operation.

**Algorithm 1** An $O(n^2)$ algorithm for Problem 1.

```
Input: $x, y \in \Sigma^+$
Output: $HCD_b(x, y)$
1: $DP[i][j] = \infty, \forall 1 \leq i \leq j \leq n$
2: Find all pairs $(i, j)$ such that $x = y[i \ldots j]$ and set $DP[i][j] = 0$.
3: for $len \leftarrow m + 1$ to $n$ do
4:   for $i \leftarrow 1$ to $n - len + 1$ do
5:      $j \leftarrow i + len - 1$
6:      if $DP[i][j] = \infty$ then
7:         $x \leftarrow \text{min-edge}_{DSR}(j - k\text{-Border}(s[i \ldots j]), j - 1)$
8:         $y \leftarrow \text{min-edge}_{DSL}(i + 1, i + k\text{-Border}(s[i \ldots j]))$
9:         $DP[i][j] = \min(x, y) + 1$
10:     end if
11:    add-leaf$_{DSR}(j, j - 1, DP[i][j])$
12:    add-leaf$_{DSL}(i, i + 1, DP[i][j])$
13:  end for
14: end for
15: return $DP[1][n]$
```

▷ **Theorem 19.** Algorithm 1 solves Problem 1 in $O(n^2)$. 
Deletion Graph

As in [18], we are interested in constructing the table $DP[i][j]$. We replace the segment tree with a linked list and change the order of processing the cells. Taking into account the iteration order (increasing according to the difference $\Delta$ of the elements in the set $R_i$), we say that it takes a step down (resp. leftward). For example, if $A_k$ is a path in $G$, we call edges from $A_k$ to a suffix $[i+\Delta]$ a downward edge and an edge from $A_k$ to a prefix $[i\ldots j-\Delta]$ a leftward edge. When a path in $G_h(x)$ uses a downward (resp. leftward) edge, we say that it takes a step down (resp. leftward). For example, if $P = (v_1, v_2, \ldots, v_s)$ is a path in $G_h(x)$, and the edge $(v_{s-1}, v_s)$ is a downward (resp. leftward) edge, we say that $P$ ends with a step left (resp. down).

Proof.

Correctness. We prove the correctness of the algorithm by induction over the algorithm’s execution. The base cases correspond to the substrings $y[i\ldots j] = x$. In these cases, $DP[i][j] = 0$ because no operation is needed to convert $x$ to $y[i\ldots j]$. Suppose we want to calculate the value of $DP[i][j]$. We remind that $DP[i][j] = \min(\min L_j, \min R_i) + 1$. We can rewrite the elements of the set $R_i$ in the following form $DP[i][p]$ with $j = \Delta(i, j) + k \leq p < j$ and the elements of the set $L_j$ in the form $DP[s][j]$ with $i < s \leq i + \Delta(i, j) - k$. Taking into account the iteration order (increasing according to the difference $j - i$) and $j > i$, we obtain the following inequalities: $j - i > j - s$ and $j - i > p - i$. Thus, it is guaranteed that when we want to calculate $DP[i][j]$ all the necessary values are already calculated.

Complexity. Line 1 runs in $O(n^2)$ and Line 2 in $O(n + m)$. For each cell above the main diagonal we have two queries and two updates both done in $O(1)$ amortized time. The overall time complexity is therefore $O(n^2)$.

4 Minimum distance common hairpin completion ancestor

In this section we study Problem 2.

Our algorithm is based on the dynamic programming technique described in [18], but we replace the segment tree with a linked list and change the order of processing the cells in the matrix. Without loss of generality, we assume that $|x| \geq |y|$ and $n = |x|, m = |y|$. As in [18], we are interested in constructing the table $DP_k[1\ldots n][1\ldots n]$ with $DP_k[i][j] = HCD_k(x[i\ldots j], \ldots)$. Similarly, we would like to compute a table $DP_k[1\ldots m][1\ldots m]$ with $DP_k[i][j] = HCD_k(y[i\ldots j], \ldots)$.

We are interested in rephrasing the problem of computing $DP_k$ in terms of shortest distances in a graph. We present the following definition.

Definition 20 (Hairpin Deletion Graph). For a string $x[1\ldots n]$, we define the Hairpin Deletion Graph $G_h(x) = (V, E)$ of $x$ as follows.

- $V = \{x[i\ldots j] | 1 \leq i \leq j \leq n\}$ is the set of substrings of $x$.
- $E = \{(x[i\ldots j], x[a\ldots b]) | x[i\ldots j] \in HR_k(x[a\ldots b])\}$ I.e. there is a directed edge from substring $A$ to the substring $B$ if $A$ can be obtained from $B$ by applying a single hairpin completion operation.

It is easy to see that $HCD_k(x[i\ldots j], x)$ is exactly the length of the shortest path from $x[1\ldots n]$ to $x[i\ldots j]$ in $G_h(x)$. Following Observation 8, we present the following characterization of the edges in $G_h(x)$.

Corollary 21. Let $x[1\ldots n]$ be a string and let $A = x[i\ldots j]$ be a substring of $x$ with $k$-Border length $l$. The set of edges emerging from $A$ in $G_h(x)$ is

$$E_A = \bigcup_{p \in [1\ldots l]} \{(A, x[i+p\ldots j]), (A, x[i\ldots j-p])\}$$

We call edges from $A = x[i\ldots j]$ to a suffix $[i+p\ldots j]$ a downward edge and an edge from $A$ to a prefix $x[i\ldots j-p]$ a leftward edge. When a path in $G_h(x)$ uses a downward (resp. leftward) edge, we say that it takes a step down (resp. leftward). For example, if $P = (v_1, v_2, \ldots, v_s)$ is a path in $G_h(x)$, and the edge $(v_{s-1}, v_s)$ is a downward (resp. leftward) edge, we say that $P$ ends with a step left (resp. down).
Then, the algorithm computes the cells of $DP_x$ row by row from top to bottom, iterating a row in decreasing order of the columns. Formally, when iterating the cell $DP_x[i][j]$, the algorithm have already computed the cells $DP_x[a][b]$ with $a < i$ and the cells $DP_x[i][b]$ with $b > j$. The order of the iteration implies a total order on the pairs $i, j \in [n] \times [n]$.

**Definition 22 (Iteration Order).** For two pairs of integers $(i_1, j_1), (i_2, j_2) \in [n] \times [n]$, we say that $(i_1, j_1)$ precedes $(i_2, j_2)$ (denoted as $(i_1, j_1) < (i_2, j_2)$) if the cell $DP_x[i_1][j_1]$ is iterated before $DP_x[i_2][j_2]$ by our algorithm. Similarly, we say $(i_2, j_2)$ proceeds $(i_1, j_1)$.

We proceed to introduce a useful concept used by the algorithm.

**Definition 23 (Restricted Path).** For a string $x[1 \ldots n]$ and integers $i, j \in [n]$, a path $P = (x, v_1, v_2, \ldots, v_z, A)$ from $x$ to $A$ in $G_h(x)$ is $(i, j)$-restricted if for every $r \in [z]$ we have $v_r = x[a_r \ldots b_r]$ such that $(a_r, b_r)$ proceeds $(i, j)$. For $(i, j) = (0, 0)$, we say that there is no $(0, 0)$-restricted path.

For integer pairs $(i, j), (a, b) \in [n][n]$ such that $(i, j) < (a, b)$, we denote as $ResL(i,j) ((a, b)]$ the length of the shortest $(i, j)$-restricted path from $x$ to $x[a \ldots b]$ in $G_h(x)$ that ends with a step left. Similarly, we denote as $ResD(i,j) ((a, b)]$ the length of the shortest $(i, j)$-restricted path from $x$ to $x[a \ldots b]$ in $G_h(x)$ that ends with a step down.

We make the following observations regarding the structure of $ResL(i,j) ((a, b)]$ and $ResD(i,j) ((a, b)]$.

**Lemma 24.** For every $i, j \in [n]$ and $a, b \in [n] \times [n - 1]$ such that $(i, j) < (a, b)$ it is satisfied that $ResL(i,j)((a,b+1]) \leq ResL(i,j)((a,b)]$.

**Proof.** If there is no $(i, j)$-restricted path that ends with a step leftwards from $(0, 0)$ to $(a, b)$, the claim is vacuously true. Otherwise, let $P = (1,n), (x_1, y_1), (x_2, y_2) \ldots (x_d, y_d), (a, b)$ be the shortest $(i, j)$-restricted path from $(1,n)$ to $(a, b)$ that ends with a step to the left. Since $P$ ends with a step leftwards, we have $x_d = a$ and there is an edge from $(a, y_d)$ to $(a, b)$. According to Corollary 21, there is also an edge from $(a, y_d)$ to $(a, b+1)$. Therefore, we can replace $(a, b)$ with $(a, b+1)$ in $P$ to obtain an $(i, j)$-restricted path $P'$ with length $|P|$ from $(1,n)$ to $(a, b+1)$.

The following symmetric statement can be proven in a similar manner.
Lemma 25. For every $i, j \in [n]$ and $a, b \in [n - 1] \times [n]$ such that $(i, j) < (a, b)$ it is satisfied that $ResD_{(i,j)}[(a, b)] \leq ResD_{(i,j)}[(a + 1, b)]$.

Lemma 24 and Lemma 25 suggest that the values of $ResL_{(i,j)}$ (resp. $ResD_{(i,j)}$) in every row (resp. column) are monotonic.

For every row $k \in [n]$ of $DP_z$, the algorithm maintains a corresponding double-sided linked list $Row_k$. Similarly, for every column $k \in [n]$ the algorithm maintains a list $Col_k$. Conceptually, $Row_k$ (resp. $Col_k$) compactly represents the values $ResL_{(i,j)}[(a,b)]$ (resp. $ResD_{(i,j)}[(a,b)]$) for all the cells $(a,b)$ in row $k$ (resp. in column $k$). Every list stores a sequence of pairs of integers $(\delta, \beta)$. The first value $\delta$ is called the distance and the second value $\beta$ is called the boundary. We call such pairs boundary pairs. The pairs are stored in increasing order of distances. For an integer $x$, we call the pair $(\delta, \beta)$ in a list the boundary predecessor (resp. boundary successor) of $x$ in $Row_k$ if $\beta$ is the minimal (resp. maximal) boundary in the list that is at least (resp. at most) $x$.

When processing $DP_z[i][j]$, we are interested in maintaining the following invariant regarding the pairs stored in $Row_a$ (for every $a \in [n]$):

Let $b \in [n]$ be an integer such that $(i, j) < (a, b)$ and let $(\delta, \beta)$ be the boundary predecessor of $b$ in $Row_a$. It holds that $ResL_{(i,j)}[(a,b)] = \delta$. Equivalently: Let $(\delta_1, \beta_1), (\delta_2, \beta_2) \ldots (\delta_z, \beta_z)$ be the pairs in $Row_a$. Note that due to Lemma 24, the pairs in $Row_b$ are naturally stored in decreasing order of their boundaries. If an integer $b$ satisfies $b \in [\beta_r \ldots \beta_{r-1} - 1]$ for some $r \in [z]$, then $ResL_{(i,j)}[(a,b)] = \delta_r$. For a visualization, see Figure 4.

Essentially, the list $Row_a$ stores an implicit representation of the shortest $(i, j)$-restricted paths that end with a left step to the cells in row $a$.

Similarly, the list $Col_b$ stores an implicit representation of $ResD_{(i,j)}[(a,b)]$ for vertices in column $b$ as follows. For every $a \in [n]$ such that $(i, j) < (a, b)$ with boundary successor $(\delta, \beta)$ in $Col_b$, it holds that $ResD_{(i,j)}[(a,b)] = \delta$.

Throughout the iterations, we maintain the pair $r = (\delta_r, \beta_r)$ such that when we iterate $DP_z[i][j]$, the pair $r$ is the boundary predecessor $j$ in $Row_i$. We also store $n$ pairs $c_1, c_2 \ldots c_n$ such that when iterating $DP_z[i][j]$, the pair $c_j = (\delta_j, \beta_j)$ is the boundary successor of $i$ in $Col_j$. We initialize every list $Row_k$ with a single pair $(\infty, 1)$ and every list $Col_k$ with a single pair $(\infty, n)$. For the sake of consistency, we treat the initialization of the algorithm as a phase in the iteration in which a dummy cell $(0, 0)$ is the currently iterated cell. Note that the initialization for the lists suggests that for every vertex, there is no $(0, 0)$-restricted path that ends with a step to downwards or leftwards.
Processing a cell. When processing \( DP_x[i][j] \), we first obtain the distance to \( DP_x[i][j] \) using \( r \) and \( c_j \). The shortest path to \((i, j)\) must end with a step to the left from a vertex in row \( i \) or with a step downwards from a vertex in column \( j \). Note that all the cells to the right of \((i, j)\) and above it have already been processed. It follows that the shortest path to \((i, j)\) is an \((i', j')\)-restricted path with \((i', j')\) being the cell processed in the previous iteration. Let \( r = (\delta_1, \beta_1) \) and \( c_j = (\delta_2, \beta_2) \). \( r \) is the boundary predecessor of \( i \) in \( \text{Row}_i \), so according to the invariant we have \( \text{ResL}_{(i', j')} = \delta_r \). Similarly, \( \text{ResD}_{(i', j')} = \delta_1^l \). We can therefore set \( DP_x[i][j] = \min(\delta_r, \delta_1^l) \).

The remaining task is to update the lists and the pointers in a manner that preserves our invariants. We make the following claims.

Claim 26. For \((i, j) \in [n] \times [n] \) and \((a, b) \in [n] \times [n] \) such that \((i, j) < (a, b)\), if an \((i, j)\)-restricted path to \((a, b)\) visits the vertex \((i, j)\) - \((i, j)\) must be the second to last vertex in the path (i.e. the next vertex is \((a, b)\).

Proof. According to Corollary 21, every edge emerging from \((i, j)\) enter a vertex \((i', j')\) such that \((i, j) < (i', j')\). The only vertex in an \((i, j)\)-restricted path that is allowed to proceed \((i, j)\) is the destination vertex. \(<

Corollary 27. Let \((i', j') \in [n] \times [n] \) and let \((i, j) \in [n] \times [n] \) be the vertex immediately following \((i', j')\) in the iteration order. Let \((a, b) \in [n] \times [n] \) such that \((i, j) < (a, b)\). If there is no edge from \((i, j)\) to \((a, b)\) we have \( \text{ResL}_{(i, j)}[(a, b)] = \text{ResL}_{(i', j')}[(a, b)] \) and \( \text{ResD}_{(i, j)}[(a, b)] = \text{ResD}_{(i', j')}[(a, b)] \)

Furthermore, by Corollary 27 and Corollary 21 together, we obtain the following.

Corollary 28. For \( k \neq i \) (resp. \( k \neq j \)), the list \( \text{Row}_k \) (resp. \( \text{Col}_k \)) does not need to be updated after the cell \((i, j)\) is processed in order to satisfy the invariant.

It follows from Corollary 28 that we only need to update the lists \( \text{Row}_i \) and \( \text{Col}_j \) to represent shortest \((i, j)\)-restricted paths instead of representing shortest \((i', j')\)-restricted paths. In other words, we need to update \( \text{Row}_i \) and \( \text{Col}_j \) to consider paths that use the vertex \((i, j)\). Specifically, paths that use \((i, j)\) as a second to last vertex (Claim 26).

We update the lists as follows. Let \( l \) be the \( k \)-Border of \( x[i \ldots j] \) and let \( d \) be the recently calculated \( d = DP_x[i][j] = \min(\delta_r, \delta_1^l) \). According to Corollary 21, there is an edge from \((i, j)\) to \((i, j - z)\) with \( z \in [1 \ldots l] \), and only to those vertices in the \( i' \)th row. We call these vertices the contested vertices. For every contested vertex, there is an \((i, j)\)-restricted path that ends with a step to the left via the vertex \((i, j)\). This path has length \( d + 1 \). Our task is to update \( \text{Row}_i \) such that every contested vertex \((a, b)\) in the list with \( \text{ResL}_{(i', j')}[(a, b)] > d + 1 \) is updated to have \( \text{ResL}_{(i, j)}[(a, b)] = d + 1 \). Every \((a, b) \in \text{Row}_i \) with \( \text{ResL}_{(i', j')}[(a, b)] \leq d + 1 \) needs to keep its current distance. The distances to uncontestd vertices in the \( i' \)th row do not require an update (Corollary 27).

Assume w.l.o.g that \( d = \delta_r \) (the case in which \( d = \delta_1^l \) is treated symmetrically). We may need to add the boundary pair \((d + 1, j - l)\) to \( \text{Row}_i \) to represent the newly available \((i, j)\)-restricted paths. First, observe that \( r = (d, \beta_r) \) should not be removed from \( \text{Row}_i \). This is due to the cost of the newly available paths via \((i, j)\) being \( d + 1 \) - longer than the paths already represented by \( \text{Row}_i \) for the vertices \((i, b)\) with \( b \in [\beta_r \ldots j] \). We follow the list pointer from \( r \) to obtain its boundary predecessor \( r' = (\delta_1, \beta_1) \) in \( \text{Row}_i \) with \( \beta_1 < \beta_r \) and \( \delta_1 > d \). We consider the following cases.
We use the algorithm required for the next iterated cell. If such that the first time, which happens of DP can be removed at most once throughout the algorithm - the overall time complexity for and to the list the process of maintaining the lists invariant, at most one pair is added to the list ▶ of Col the new boundary predecessor (resp. boundary successor) of the next pair in the list until a pair execute the required deletion in a straightforward manner. Starting from represents a path with length at least may be insufficient. We also need to remove every pair (δ, β) in Rowi with β ∈ [j − l...βl]. All of those pairs are now redundant in Rowi - as they represent paths with a length at least d + 1 to contested vertices. We execute the deletion of these pairs in a straightforward manner by following the links from r′ until we reach a pair r∗ = (δ, β) with β < j − l. When r∗ is finally met, we insert (d + 1, j − l) to Rowi between the r and r∗.

We proceed to treat Colj. If δj = d, the treatment of Colj is completely symmetric to the treatment of Rowi. Otherwise, δj. As in the treatment of Rowi, our task is to add a representation of the paths with length d + 1 to vertices (a, j) with a ∈ [i...i + l]. Namely, every pair (δ, β) in Colj with δ < i + l should be removed (including c1), as it represents a path with length at least d + 1 to one of the vertices (a, j) with a ∈ [i, i + l]. We execute the required deletion in a straightforward manner. Starting from c1, we proceed to the next pair in the list until a pair (δ, β) with β > j + l is found. We then remove the pairs iterated in this process from Colj and append (d + 1, i + l) to the beginning of the list. This concludes the updates to Rowi and to Colj. We note that if r of cj is removed from Rowi or from Colj, respectively, the new pair (d + 1, j − l) (or respectively, (d + 1, i + l)) is becoming the new boundary predecessor (resp. boundary successor) of j in Rowi (resp. of i in Colj).

Finally, we need to update r and cj to be the boundary predecessor and successors required for the next iterated cell. If i < n, the i value will remain the same on the next iteration. In this case, if βr = j, we update r to be the next element in Rowi. If βr < j, we do not need to update r is it is also the boundary predecessor of j − 1 in Rowi.

If i = n, the next iteration is the first step in row i + 1. It follows that Rowi+1 is still in its initialized state, and we set the only pair in (∞, 1) ∈ Rowi+1 to be r.

As for the cj pointers, we need to update all of them every time a new row is met. When moving from row i to row i + 1, every cj needs to be updated from the predecessor of i in Colj to the predecessor of i + 1 in Colj. This is done in a symmetric manner to the update of r.

Case 1.a: δ1 = d + 1 and j − l ≥ β1. In this case, Rowi already represents the shortest restricted paths with cost d + 1 to the vertices (i, k) with k ∈ [j − l...βl − 1]. Therefore, no update is required for Rowi.

Case 1.b: δ1 > d + 1 and j − l ≥ β1. In this case, we need to add the boundary pair (d + 1, j − l) after the boundary border r in Rowi. The following pairs in Rowi have boundaries smaller than j − l and therefore represent the shortest paths uncontested vertices and do not need to be changed. If j − l = β1, we also remove r′ from Rowi, as it is redundant.

Case 2 : j − l < β1. In this case, adding the pair (d + 1, j − l) after the pair r to Rowi may be insufficient. We also need to remove every pair (δ, β) in Rowi with β ∈ [j − l...βl]. For the treatment of MDCA, this can return the answer in CPM 2023.
Figure 5 Case 2: The shortest path to the purple vertex is discovered to be 2. The $k$-Border of the substring corresponding to this vertex is 11 (denoted as the purple left arrow), thus creating new restricted paths to the vertices covered by the purple arrow (recall that in $G_h(x)$, there is a directed edge from the purple vertex to every one of the vertices covered by the purple arrow). The appropriate update to the list is removing the pairs $(3,10)$ and $(4,6)$ as the vertices in the green area and in the blue area are now accessible via a shorter path with length 2 via the purple vertex. This new paths are represented by the newly added pair $(2,5)$.

Figure 6 Case 1.a: The distance to the purple vertex is discovered to be 3, enabling new paths with length 4 to the vertices touched by the purple arrow (representing the length of the $k$-Border of the substring corresponding to the purple vertex). These new paths do not improve upon the restricted paths already represented in the list, so the pair $(4,8)$ representing these new paths is simply not added to the list.

$O(n^2)$. To be clear, we provide a brief explanation of the algorithm. In the first stage, the algorithm builds a trie with all the suffixes of the string $x$. Then, it will traverse the trie for every suffix of $y$ and at every match it will compute the sum of $DP$ values. In short, this algorithm determines in quadratic time all the common substrings of $x$ and $y$ and keeps the one with the minimum sum of distances. We add the pseudocode for the algorithm described in this section in the appendix.
Figure 7 Case 2: The distance to the purple vertex is discovered to be 3. This creates new restricted paths with length 4 to the vertices touched by the purple arrow (representing the k-Border of the string corresponding to the purple vertex). For the vertices in the green area, this is not an improvement, as we already have a representation to a path with length 3 to those vertices. The distances to the vertices in the blue area and to the vertex in the red area touched by the purple arrow are longer or equal to 4. To represent this, we add the boundary pair (4, 5) and remove the boundary pair (4, 6) (as the k-Border (4, 6) represented the distances to the vertices in the blue interval, which are now represented by (4, 5).

5 Conclusions and future work

In this paper we study two problems related to the hairpin completion operation. We propose a quadratic time algorithm for solving these two problems, thus improving the runtime over previous work by Manea [18]. Notice that both our algorithms compute the dynamic programming table of the respective problem explicitly.

A question that arises from our work is can one find an algorithm that solves one of these problems by computing a small subset of cells in the dynamic programming table, which implies a runtime of \( o(n^2) \). An interesting and challenging open problem is to provide an \( o(n^2) \) algorithm for any of the two problems studied in this paper (not necessary with uses of the dynamic programming’s formula), or present a lower bound matching with known problems.

For other variants of hairpin problems (see, e.g., [9, 20, 21]), we believe our techniques can help understand them better and help with designing efficient algorithms for these problems.

References


### Appendix

Algorithm 2

An $O(n^2)$ algorithm for Problem 2.

**Input:** $x, y \in \Sigma^+$

**Output:** a string $z$ such that $HCD_k(z, x) + HCD_k(z, y)$ is minimum

1. $DP_x = \text{Compute}DP(x)$
2. $DP_y = \text{Compute}DP(y)$
3. return $\text{Compute}_MDCA(x, y, DP_x, DP_y)$

Algorithm 3

Updates the list $Row_i$.

1. procedure $\text{UPDATE}ROW(i, j, r)$
2. while $r$ is not NULL and $\beta_r > j - k$-Border($s[i...j]$) and $\delta_r > DP_x[i][j]$ do
3. delete $r$ from $Row_i$
4. $r = r \rightarrow \text{next}$
5. end while
6. add $(j - k$-Border($s[i...j]$), $DP_x[i][j] + 1)$ to $Row_i$
7. end procedure

Algorithm 4

Updates the list $Col_j$.

1. procedure $\text{UPDATE}COL(i, j, c_j)$
2. while $c$ is not NULL and $\beta^l_c < i + k$-Border($s[i...j]$) and $\delta^l_c > DP_x[i][j]$ do
3. delete $c_j$ from $Col_j$
4. $c_j = c_j \rightarrow \text{next}$
5. end while
6. add $(i + k$-Border($s[i...j]$), $DP_x[i][j] + 1)$ to $Col_j$
7. end procedure
Algorithm 5 ComputeDP.

Input: $x \in \Sigma^+$
Output: $DP_x$

1. $DP[i][j] = \infty, \forall 1 \leq i \leq j \leq n$ \hspace{1cm} ▷ $n$ is the length of the input string
2. $DP_x[1][n] = 0$ \hspace{1cm} ▷ Base case
3. add $(n - k$-Border$(s[1 \ldots n]), 1)$ to Row$_1$ \hspace{1cm} ▷ Compute the first line of $DP_x$
4. for $i \leftarrow n - 1$ to 1 do
5. \hspace{1cm} if $i \leq \beta_r$ then
6. \hspace{2cm} $DP_x[1][i] = \delta_r$
7. \hspace{1cm} if $j - k$-Border$(s[1 \ldots i]) < \beta_r$ then
8. \hspace{2cm} updateRow$(1, i, r)$
9. \hspace{1cm} end if
10. \hspace{1cm} end if
11. \hspace{1cm} end for
12. for $i \leftarrow 2$ to $n$ do
13. \hspace{1cm} for $j \leftarrow n$ to 1 do
14. \hspace{2cm} if $\delta_r < \delta^j_i$ then
15. \hspace{3cm} if $j \geq \beta_r$ then
16. \hspace{4cm} $DP_x[i][j] = \delta_r$
17. \hspace{3cm} if $j - k$-Border$(s[i \ldots j]) < \beta_r$ then
18. \hspace{4cm} updateRow$(i, j, r)$
19. \hspace{4cm} updateCol$(i, j, c_j)$
20. \hspace{3cm} end if
21. \hspace{3cm} end if
22. \hspace{2cm} else
23. \hspace{3cm} if $i \leq \beta^j_i$ then
24. \hspace{4cm} $DP_x[i][j] = \delta^j_i$
25. \hspace{3cm} if $i + k$-Border$(s[i \ldots j]) > \beta^j_i$ then
26. \hspace{4cm} updateCol$(i, j, c_j)$
27. \hspace{4cm} updateRow$(i, j, r)$
28. \hspace{3cm} end if
29. \hspace{3cm} end if
30. \hspace{2cm} end if
31. \hspace{1cm} end for
32. \hspace{1cm} end for
33. return $DP_x$
On Distances Between Words with Parameters

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Abstract

The edit distance between parameterized words is a generalization of the classical edit distance where it is allowed to map particular letters of the first word, called parameters, to parameters of the second word before computing the distance. This problem has been introduced in particular for detection of code duplication, and the notion of words with parameters has also been used with different semantics in other fields. The complexity of several variants of edit distances between parameterized words has been studied, however, the complexity of the most natural one, the Levenshtein distance, remained open.

In this paper, we solve this open question and close the exhaustive analysis of all cases of parameterized word matching and function matching, showing that these problems are \(\mathsf{NP}\)-complete. To this aim, we also provide a comparison of the different problems, exhibiting several equivalences between them. We also provide and implement a MaxSAT encoding of the problem, as well as a simple FPT algorithm in the alphabet size, and study their efficiency on real data in the context of theater play structure comparison.

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1 Introduction

Measuring the similarity between text strings is a fundamental problem in computer science, and has applications in bioinformatics [23], databases [1, 16] and natural language processing [27]. Among the measures of similarities between strings, the Levenshtein distance [28] is the most commonly used, both for its practicality and its ease of computation. This distance quantifies the minimum number of operations of insertion, deletion, and substitution needed to transform a string into another one. It has a wide range of applications, ranging from biological sequence alignment [33] to dialect pronunciation differences [25] or signature authentication [34]. Computing the edit distance between two strings of length \(n\) and \(m\) can be achieved in time \(O(nm)\), by computing the distance between all their prefixes, and storing the results in a dynamic programming fashion [37]. The success of the Levenshtein distance generated many extensions and generalization on more complex models, such as trees [38] or automata [32].

However, a limitation of the Levenshtein distance is that it only captures proximity between strings (or objects) written on the same alphabet. Evaluating the proximity of strings written on different alphabets is a problem that arises in various applications, such as bioinformatics [35], image processing [17] and code duplication [6, 7]. In all those contexts,
the technique used is the one of parameterized matching [6, 7]. Instead of using classical strings, parameterized matching uses “parameterized words” written using both constant parts, which are expensive to rename, and parameters, which are meant to be renamed freely. Formally, two equal-length strings $u$ and $v$ over an alphabet $\Pi$ are said to be parameterized matching if there exists a 1-to-1 function $f : \Pi \to \Pi$ such that $f(u) = v$, where $f(u)$ is defined as $f(u_1) \ldots f(u_{|u|})$.

Words with parameters also occur in other frameworks, and are often used in slightly different ways. The first of those frameworks was initially introduced in the context of Ramsey theory in the 80s [36], and is called “parameter words”. In this context, parameters are labelled according to their order of first occurrence. Parameter words are also equipped with a composition operation, where parameters of the first word can be instantiated by characters or parameters of the second word. Parameter words can also be seen as equivalence classes of parameterized words, which are the main focus of this article.

A second framework using parameters is the one of parameterized regular expressions introduced in [10], where parameters can only be instantiated by constants, and not by other parameters. The focus in this context is therefore made on the set of all possible valuations of the parameters. Then, when defining algorithmic problems on such objects, two distinct semantics can be studied: either the “certainty semantics”, where all valuations need to have some property, or the “possibility semantics”, where at least one valuation needs to have this property. To make a difference with the parameterized word framework mentioned below, we choose to call these words “instantiable words”. Finally, this notion of words with parameters can also be seen as a refined version of partial words (words containing a wildcard character) [15]. The notion of partial words is also important in the context of databases where paths of incomplete graphs can be interpreted as instantiable words [9].

This article aims at studying similarity by using edit distances in the framework of words with parameters. In this framework, the pattern matching problem, which consists in looking for the first string as a subword of the second string, has been extensively studied, either looking for exact occurrences, with efficient algorithms [4, 19, 30] or approximate ones, which is often NP-hard [21, 22]. In the case where we compare the two input strings in their entirety, various exact parameterized matching problems have been studied for parameterized pattern matching, namely string parameterized matching [7], single pattern parameterized matching [7, 3], multiple pattern parameterized matching, or 2-dimensional parameterized matching, many of those works being compiled in [29] and [31]. Different approximate variants of parameterized matching using edit distance have already been studied, but the problem has not been completely solved: the first work on the topic is [8], in which Baker introduces a form of approximate parameterized pattern matching in which the replacement of any substring by another one that is in parameterized matching with it is considered as a base edit operation. Parameterized matching under the Hamming distance, i.e., with a distance allowing only substitutions, has been covered in [24], where the authors prove that the string matching problem with at most $k$ mismatches can be solved in time $O(m + k^{1.5})$. The LCPS (Longest Common Parameterized Subsequence) problem, equivalent to the parameterized pattern matching problem with insertions and deletions, is shown to be NP-hard in [26], which also provides an approximation algorithm. Those two different complexity classes for these problems raise the question of the complexity of the problem under the Levenshtein distance. This problem was left as an open question in the conclusion of [24].

Our paper establishes that this problem is NP-complete. Moreover, the result also extends to any possible edit distances obtained from deletion, insertion, and substitution as soon as substitution is not the only operation allowed, as summarized in Figure 1. Our main
proof also implies the main theorem of [26] with a different \( \text{NP} \)-completeness reduction. This contrasts with the problems of exact parameterized pattern matching which are all polynomial-time solvable, as well as all variants of the string matching problem with deletions, insertions or substitutions.

We also extend these results to function matching, which is the problem obtained by relaxing the 1-to-1 restriction in parameterized matching, as defined in [2]. This generalization, by breaking the symmetry of parameterized matching, actually gives rise to two close but different problems, depending of the order of operations that are considered. We study the links between all these problems and their computational complexity, and study two practical ways to solve them, parameterized complexity and the use of maxSAT solvers.

We also make a direct connection with the framework of instantiable words, more precisely with a natural problem of distance between languages. We show how instantiable word problems can be reduced to parameterized matching ones, under the right assumptions. This allows us to open new perspectives on the complexity of several language repair problems.

In Section 2, we give basic definitions and notations, and recall the existing formalism of parameterized matching and instantiable words. In Section 3 we discuss approximate parameterized matching and its various generalizations. We also link it to instantiable words. In Section 4, we first prove a collection of technical results that build up to the \( \text{NP} \)-completeness proofs for parameterized matching and function matching problems defined above. In Section 5, we study two approaches to solve one of the variants of parameterized matching in practice, a simple FPT algorithm parameterized by the alphabet size and a MaxSAT encoding. We show in Section 6 that these implementations can solve real instances of the problem, motivated by structure comparison of theater plays.

Finally, in Section 7, we conclude the paper and give a few perspectives on the notion of distance between parameterized languages.

## 2 Notations and Definitions

### 2.1 Basic Notations on Words and Editions

#### Words

An alphabet is a set of letters. A word on an alphabet \( A \) is a finite sequence of letters from \( A \), indexed starting from 1. Let \( u \) be a word on \( A \). Unless defined differently, we note \( u_i \) the \( i \)-th letter of \( u \), and \( |u| \) is the length of \( u \). If \( i \not\in [1, |u|] \), \( u_i \) is defined as the empty word \( \varepsilon \). If \( x \) is a letter from \( A \), \( |u|_x \) is the number of times \( x \) appears in \( u \). Similarly, if \( X \) is a set of letters, \( |u|_X = \sum_{x \in X} |u|_x \) is the number of occurrences of letters of \( X \) in \( u \). If \( f \) is a function defined on an alphabet \( A \), we extend it to \( A^* \) in the usual way, so that \( f(u) = f(u_1) \ldots f(u_{|u|}) \).

If \( f \) is a function, we denote by \( \mathcal{D}(f) \) the domain of \( f \). Two functions \( f \) and \( g \) are said to be compatible if \( f|_{\mathcal{D}(g) \cap \mathcal{D}(f)} = g|_{\mathcal{D}(g) \cap \mathcal{D}(f)} \). The identity function on \( D \) is defined as \( \text{Id}_D(x) = x \) for all \( x \) in \( D \).

<table>
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<tr>
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<th>I \footnote{Cor. 14}</th>
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<tr>
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<td>P \footnote{[24]}</td>
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<td>\text{NP}\text{-}\text{complete} (Cor. 14)</td>
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**Figure 1** Complexity of the variants of parameterized matching \( PM^d \), depending on the kind of operations (D: deletion, I: insertion, S: substitution) allowed in the edit distance \( d \).
6.4 On Distances Between Words with Parameters

**Edit Operations**

In this paper, we consider the three classical edit operations which are deletion, substitution and insertion. Let \( u = u_1 \ldots u_n \) be a word of size \( n \). Let \( i \) be an integer between 0 and \( n \) and \( x \) be a letter of the alphabet, the insertion at position \( i \) is the operation that transforms \( u \) to \( u_1 \ldots xu_{i+1} \ldots u_n \). Let \( j \) be an integer between 1 and \( n \), the deletion at position \( j \) is the operation that transforms \( u \) into \( u_1 \ldots u_{j-1}u_{j+1} \ldots u_n \). Let \( y \) be a letter of the alphabet and \( y \neq u_j \), the substitution to \( y \) at position \( j \) is the operation that transforms \( u \) into \( u_1 \ldots u_{j-1}yu_{j+1} \ldots u_n \). A sequence of operations or rewriting sequence \( \rho \) is a sequence of edit operations. We denote by \( \rho(u) \) the word obtained by applying the edit operations of \( \rho \) one after another, in the order defined by \( \rho \), to \( u \).

**Distances**

Given a set of edit operations \( E \) and two words \( u \) and \( v \), the edit distance between \( u \) and \( v \) is defined as the length of a shortest sequence of operations of \( E \) changing \( u \) into \( v \). We denote by \( D \) the distance obtained on words by allowing only deletion operations: that is to say \( D(u,v) = k \) iff \( v \) can be obtained by deleting \( k \) letters from \( u \). Similarly, we note \( I \) and \( S \) the distances obtained by allowing only insertions and substitutions respectively (note that \( S \) is the Hamming distance). We also combine these notations to define \( DI \) as the distance with insertions and deletions, and so on. We also denote the Levenshtein distance \( DIS \) by \( L \). Note that some of these edit distances are not metrics, because they are not symmetrical. We emphasize this by calling symmetric edit distances the distances \( DI, S, \) and \( L \).

2.2 Comparing Words with Parameters

Conceptually, a word with parameters is a word in which some letters are not yet determined. In order to distinguish the parameters from the constants, we split the alphabet into \( \Sigma \), the alphabet of the constants and \( \Pi \), the alphabet of the parameters. By definition, these alphabets are finite. A word with parameters can either be seen as representing a “word template” (i.e., a word with variable parts), or a set of words (determined by all possible affectations of its parameters). Depending on the definition chosen, comparing two words \( w_1 \) and \( w_2 \) is done in two different ways. In the first setting [6, 7, 8, 31, 2, 5, 24, 29, 26, 17], parameters of \( w_1 \) are renamed through a function \( f \) that maps the set of parameters to itself, and acts as identity on the set of constants. It is then possible to compare \( f(w_1) \) and \( w_2 \), which are written on the same alphabet. In the second setting, constants are seen as the concrete values parameters can take [11]. Parameters are instantiated through two functions \( f_1 \) and \( f_2 \) that map constants to themselves, but also map parameters to constants. The words \( f_1(w_1) \) and \( f_2(w_2) \) are then made only of constants, and can be compared. Formally, this gives rise to the two following different definitions:

On the one hand, a parameterized word is a word on an alphabet \( \Sigma \cup \Pi \). In all that follows, \( \Sigma \) and \( \Pi \) are two disjoint alphabets, respectively called the alphabet of constants and the alphabet of parameters. Alphabets \( \Sigma \) and \( \Pi \) are considered to be finite, unless specified otherwise.

Two parameterized words \( u \) and \( v \) are said to be in function matching if there exist \( f_1 : \Pi \to \Pi \) and \( f : \Pi \cup \Sigma \to \Pi \cup \Sigma \) such that \( f|_\Pi = f_1 \), \( f|_\Sigma = Id_\Sigma \), and \( f(u) = v \). In the classical setting [6], \( f \) is also constrained to be 1-to-1, and this relationship is called parameterized matching. Note that parameterized matching is an equivalence relation on parameterized words. Testing if two words are parameterized matching can be achieved in linear time [7].
On the other hand, an instantiable word is a word on the alphabet $\Sigma \cup \Pi$. Given $f : \Pi \to \Sigma$, we extend it to constants by setting $f(x) = x$ for all $x \in \Sigma$, and we then define the language of an instantiable word $u$ to be $L(u) = \{w \in \Sigma^* \mid \exists f : \Pi \to \Sigma, f(u) = w\}$. This definition is akin to the $L\diamond$ semantic of a parameterized regular expression defined in [11], but restricted here to a single instantiable word. Two instantiable words $w_1$ and $w_2$ describe the same elements if their languages are equal, i.e. $L(w_1) = L(w_2)$.

3 Different Definitions for Different Semantics and Problems

In this section, we introduce various new approximate variants of parameterized matching, and compare them, highlighting their differences on examples.

3.1 Variants of Parameterized Matching

In parameterized matching, the function $f$ renaming parameters is generally considered to be 1-to-1. In this paper, we also consider the function matching problem, which is the case where $f$ is not constrained to be injective anymore, as defined in [2]. We also introduce multiple approximate variants of the parameterized matching problems, depending on several edit distances obtained by combining insertion, deletion and substitution operations.

3.1.1 Edit distances for parameterized matching between two strings: $PM^d$

- **Definition 1.** If $d$ is an edit distance, we denote by $PM^d$ the parameterized matching problem under $d$, which is the following:
  - **Input:** two parameterized words $u, v$, a parameter alphabet $\Pi$, an alphabet $\Sigma$ of constants, and a natural number $k$.
  - **Problem:** Does there exist $u'$ such that $d(u, u') \leq k$ and $u'$ and $v$ are parameterized matching, i.e. there exists a 1-to-1 function $f : \Pi \cup \Sigma \to \Pi \cup \Sigma$ such that $f|\Sigma = Id_{\Sigma}, f(\Pi) = \Pi$, and $f(u') = v$ ?

In that case, we say that $u'$ and $f$ realize the matching between $u$ and $v$. We sometimes write that only $f$ or $u'$ realize the matching if the other one is not relevant to a proof.

In cases where $\Sigma$ and $\Pi$ are already defined, we omit them and simply call $PM^d(u, v, k)$ the result of the decision problem. Furthermore, $PM^d(u, v)$ denotes the minimum integer $k$ (potentially infinite) such that $PM^d(u, v, k)$ is true.

We can note that this problem can be solved in polynomial time adapting the classical dynamic programming algorithm [33, 37] when the alphabet sizes are fixed.

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Figure 2 Side-by-side comparison of $PM^{DIS}$, $FM_1^{DIS}$ and $FM_2^{DIS}$.
3.1.2 Edit distances for function matching between 2 strings: $FM_1^d$

To denote function matching problems, we use $FM$ instead of $PM$: $FM_D$ denotes the function matching problems with deletions.

Furthermore, if $P$ is one of the problems defined above, we note $P_1$ the problem where edit operations are only applied to the first argument, and $P_2$ the one where they are only applied to the second argument.

**Definition 2.** The $FM^d_1$ and $FM^d_2$ problems are defined as follows. For both problems, the input is the following:

- **Input:** two parameterized words $u,v$, a parameter alphabet $\Pi$, a constant alphabet $\Sigma$, and a natural number $k$.

The problems are then:

- **Problem $FM^d_1$:** $\exists u'$ such that $d(u,u') \leq k$ and $u'$ and $v$ are function matching?
- **Problem $FM^d_2$:** $\exists v'$ such that $d(v,v') \leq k$ and $u$ and $v'$ are in function matching?

Note that the renaming function $f$ is always applied to one input only. These definitions are illustrated on an example in Figure 2.

3.2 Comparing Variants of $PM$

In this subsection, we compare the different variants of our problem.

Regarding the one-to-one parameterized matching $PM$, note that the definition we give above is designed to be easily extended to the different variants when we drop the one-to-one restriction. In [24], the authors consider that the “correct way for defining the edit distance problem” is “to allow the operations and then apply the edit distance”. By one-to-one restriction. In [24], the authors consider that the “correct way for defining the edit distance problem” is “to allow the operations and then apply the edit distance”. By extending the definition of $FM_1^d$ and $FM_2^d$ to define $PM_1^d$ and $PM_2^d$ in the case of one-to-one matching, we see that it is actually possible to switch the order of operations, and to reverse them (deletions then become insertions and vice versa, and the renaming function $f^{-1}$ is well-defined), in this case. This makes our definition consistent with the quote from [24] above. Formally, this gives the following equalities, for all parameterized words $u$ and $v$:

$$PM_1^d(u,v) = PM_D^D(v,u) = PM_D^D(u,v) = PM_1^d(v,u).$$

More generally, it holds that for every edit distance $d$, $PM_1^d(u,v) = PM_2^d(v,u) = PM_2^d(u,v)$, where $d^{-1}$ denotes the converse distance of $d$, i.e. $d^{-1}$ contains deletions if $d$ contains insertions, insertions if $d$ contains deletions, and substitutions if $d$ contains substitutions.

However, for function matching, we only have the following equalities: $FM_1^d(u,v) = FM_D^D(u,v)$ and $FM_D^D(u,v) = FM_1^d(u,v)$.

By taking $u = ab$ and $v = cc$, we can notice that $FM_1^d(u,v) = 0$ and $FM_D^D(v,u) = \infty$, so the equality $FM_1^d(u,v) = FM_D^D(v,u)$ does not hold.

Finally, note the following inequalities:

**Proposition 3.** Let $u$ and $v$ be parameterized words over $\Sigma \cup \Pi$. Then:

1. $FM_1^d(u,v) \leq PM_1^d(u,v)$;
2. If $d$ is a symmetric edit distance, $FM_1^d(u,v) \leq FM_1^d(u,v)$.

**Proof.** The first point comes from the fact that any solution to $PM_1^d$ is also a solution to $FM_1^d$. For the second point, let $k = FM_1^d(u,v)$, and let $u'$ and $f$ realize $FM_1^d(u,v)$. We construct a word $v'$, obtained by applying to $v$ the same operations applied to $u$ to obtain $u'$, but “mirrored”. That is to say, for every operation used in $u$, we apply an operation in $v$, in the following way:
If a letter $a$ is inserted in $u$, there exists a position $i$ in $u'$ such that $u'_i = a$, and $f(u'_i) = v_i$. Hence, we delete $v_i$ in $v$.

Similarly, if a letter is substituted for another letter $a'$ in $u$, there exists $i$ such that $u'_i = a'$, and we substitute $v_i$ to $f(a)$.

If a letter $a$ is deleted in $u$ at position $i$, we insert $f(a)$ in $v$ at position $i$ instead.

It then holds that $f(u) = u'$, and hence $FM^d(u, v) \leq k$.

Note that the above proof does not work to prove the converse inequality between $FM^d_1$ and $FM^d_2$, as it would require to consider an element of $f^{-1}(a)$, which might be empty. This is illustrated in the following example, on the alphabet $\Pi = \{a, b\}$:

**Example 4.** Let $N \in \mathbb{N}$ and consider $u = a^N b^N b$ and $v = a^N a^N b$. $u$ and $v$ are not in parameterized matching, hence $FM^d_{DIS}(u, v) > 0$ and $FM^d_{DIS}(u, v) > 0$. By substituting the last $b$ in $v$ for a $a$, and picking a function $f$ such that $f(a) = f(b) = a$, we get $FM^d_{DIS}(u, v) = 1$ (see Figure 2 for an example with $N = 2$). For $FM^d_{DIS}$, since $b$ appears in $v$, it holds that for any function $f$ realizing $FM^d_{DIS}$, $f(a) = b$ or $f(b) = b$. Hence, at least $N$ occurrences of $b$ appear in $f(u)$. Since there is only one occurrence of $b$ in $v$, it is clear that $FM^d_{DIS}(u, v) \geq N - 1$.

The difference between $FM^d_1$ and $FM^d_2$ comes from the fact that $\Pi$ is fixed in the input. In the case where $\Pi$ could be extended, both problems can be shown equivalent (for example if we allow a new letter $c$ in the example of Figure 2, we also get $FM^d_{DIS}(u, v) = 1$ by setting $u_5 \rightarrow c$ and $f : [a \rightarrow a, b \rightarrow a, c \rightarrow b]$), by using the same proof as Proposition 3.

### 3.3 Instantiable Words versus Parameterized Words

The parameterized word formalism and the instantiable word formalism give rise to two different definitions of distances between words. Given an edit distance $d$ on words, there are two ways to extend it to words with parameters. Let $w_1$ and $w_2$ be two words over $\Sigma \cup \Pi$. The two possible extensions are the following:

- The distance between $w_1$ and $w_2$ is defined as $d(w_1, w_2) = PM^d(w_1, w_2)$. Alternatively, the function distance between $w_1$ and $w_2$ is defined as $FM^d_1(w_1, w_2)$.

- The function distance between $w_1$ and $w_2$ is the distance between their respective languages seen as sets, that is to say $d(w_1, w_2) = d(L(w_1), L(w_2)) = \sup_{u \in L(w_1)} \inf_{v \in L(w_2)} d(u, v)$. Equivalently, $d(w_1, w_2) \leq k$ if and only if for all $f_1 : \Pi \rightarrow \Sigma$, there exists $f_2 : \Pi \rightarrow \Sigma$ such that $d(f_1(w_1), f_2(w_2)) \leq k$.

This second definition stems from the definition of distance between languages, as defined and studied in [12, 13, 14].

**Example 5.** Consider the words $u = axyb$ and $v = xbyy$, on the alphabets $\Sigma = \{a, b\}$ and $\Pi = \{x, y\}$, and consider the distance $S$. On the one hand, $FM^d_2(u, v) = 4$, because regardless of the matching chosen, every letter of $f(u)$ has to be substituted. On the other hand, for any function $f_1 : \Pi \rightarrow \Sigma$, choosing $f_2$ such that $f_2(x) = a$ and $f_2(y) = b$ yields a distance $d(f_1(u), f_2(v))$ of at most 2, by substituting the 2 middle letters.

Given a big enough alphabet, those two definitions can in fact be shown equivalent:

**Proposition 6.** Let $w_1$ and $w_2$ be words over $\Sigma \cup \Pi$, and let $d$ be a symmetric edit distance on $\Sigma \cup \Pi$. Suppose $|\Sigma| \geq |w_1| + |w_2|$, and let $k$ be an integer. Then, the following are equivalent:

1. $FM^d_1(w_2, w_1) \leq k$
2. $d(L(w_1), L(w_2)) \leq k$
Notice how \( w_1 \) and \( w_2 \) change position between the two distances. This is not benign, as \( FM^d_I \) is not symmetric.

**Proof.** Suppose \( FM^d_I(w_2, w_1) \leq k \). There exists \( f : \Pi \to \Pi \) such that \( d(f(w_2), w_1) \leq k \). For this proof, we will use the characterization of the distance between languages with \( f_1 \) and \( f_2 \). Let \( f_1 : \Pi \to \Sigma \). Define \( f_2 = f_1 \circ f \). Since \( d(w_1, f(w_2)) \leq k \), we have \( d(f_1(w_1), f_1(f(w_2))) \leq k \), by following the same edit operations. Hence \( d(f_1(w_1), f_2(w_2)) \leq k \).

Suppose now \( d(L(w_1), L(w_2)) \leq k \). Let \( f_1 : \Pi \to \Sigma \) be a 1-to-1 function such that for all parameters \( x \) in \( w_1 \), \( f(x) \) does not appear in \( w_1 \) or \( w_2 \). This is possible since \( \Sigma \) is large enough. There exists \( f_2 : \Pi \to \Sigma \) such that \( d(f_1(w_1), f_2(w_2)) \leq k \). Let \( h : \Sigma \to \Pi \cup \Sigma \) be such that if \( x \in \Pi \), \( h(f_1(x)) = x \), and if \( x \not\in f_1(\Pi) \), \( h(x) = x \). We then have \( h \circ f_1 = Id \). What is more, since \( h \) is injective, \( d(f_1(w_1), f_2(w_2)) = d(h(f_1(w_1)), h(f_2(w_2))) = d(h(f_2(w_2)), w_1) \).

Hence, \( FM^d_I(w_2, w_1) \leq k \).

### 4 Hardness Results for Approximate Parameterized Matching

In this section, we study the complexity of the various parameterized matching problems. We show the \( \mathsf{NP} \)-completeness of the simplest problems using only deletions, which will be sufficient to show the \( \mathsf{NP} \)-completeness of all the other problems. We start by proving some practical lemmas, and then proceed to the reductions.

#### 4.1 “Block by block” Lemmas

In this section, we regroup a few useful technical lemmas. We start of by stating two simple results on distance and words, for which the proofs can be found in Appendix A. We then turn to block lemmas, which will later be useful in the proofs of Theorems 12,17 and 15, to combine the various gadgets defined during the reduction.

This lemma simply states a commutativity result between the application of a matching \( f \) and the rewriting steps.

**Lemma 7.** Let \( d \) be a distance, \( k \) an integer and \( u,v \) two parameterized words such that \( PM^d(u,v) \leq k \), and let \( f \) realize this parameterized match. Then: \( d(f(u), v) \leq k \). The same result holds for \( FM^d_I(u,v) \).

**Proof Idea.** The proof is done by induction on \( k \). We discuss whether the \( (k+1) \)-th operation is an insertion, a deletion, or a substitution, and show that a corresponding operation can be used in \( f(u) \).

**Lemma 8.** Let \( z,u \) and \( v \) be (parameterized) words, and let \( d \) be a distance. Then \( d(zu,zv) = d(u,v) \).

**Proof Idea.** We show that we can consider every rewriting operation to be applied in \( u \) only: if \( z \) is modified during an optimal rewriting sequence, the words have some redundancy, and the same operations could have been carried in \( u \) instead. We proceed again by induction, and focus on the base case by studying the 3 possible cases, one for each type of operation.

Next, we turn to prove “block by block” matching lemmas. Those results state that it is possible to encode multiple parameterized matching instances into a single one. They hold for every type of problems considered here, but their proofs vary slightly; we present them in order of increasing complexity. Note that all the constructions given can be achieved in polynomial time.
Lemma 9. Let \( u_1, \ldots, u_n \) and \( v_1, \ldots, v_n \) be parameterized words over \( \Sigma \cup \Pi \) such that for \( 1 \leq i \leq n \), \( k_i = |u_i| - |v_i| \geq 0 \), and \( k = \sum_{i=1}^{n} k_i \). There exist \( u \) and \( v \) two parameterized words over \( \{\#\} \cup \Sigma \cup \Pi \), where \( \# \) is a fresh variable, such that the following are equivalent:

1. \( PM^D(u, v) = k \)
2. For all \( 1 \leq i \leq n \), \( PM^D(u_i, v_i) = k_i \) and the applications \( f_i \) realizing those matchings are all compatible.

Proof. The idea behind this proof and all the following ones is that we can introduce a separator \( \# \) to concatenate all the words, and that this separator will never be touched by any deletions or applications of \( f \).

Let \( \# \) be a fresh constant. We define \( u = u_1 \# u_2 \# \ldots \# u_n \) and \( v = v_1 \# v_2 \# \ldots \# v_n \).

1. \( PM^D(u, v) \leq k \). Let \( u' \) and \( f \) realize this parameterized match. Since the \( \# \) symbols are constants, we have \( f(\#) = \# \). Since \( u' \) is obtained from \( u \) by deletions, we have \( |u'| |\#| \leq |u| |\#| \). Since \( f \) is injective and \( f(\#) = \# \), \( |f(u')| |\#| \leq |f(u)| |\#| \). Hence, it holds that \( |v| |\#| = |f(u')| |\#| \leq |f(u)| |\#| = |u| |\#| \). Since \( |u| |\#| = |v| |\#| \), this is an equality, and \( |f(u')| |\#| = |f(u)| |\#| \). Hence \( |u'| |\#| = |u| |\#| \), and no \( \# \) character is deleted. The word \( u' \) is then of the form \( u'_1 \# u'_2 \# \ldots \# u'_n \), where \( |u'_i| |\#| = 0 \) and \( D(u_i, u'_i) = k_i \) for all \( i \). Thus, \( f(u') = f(u'_1) f(u'_2) \ldots f(u'_n) = v_1 \# v_2 \# \ldots \# v_n \). Since no other \( \# \) letter appear in any \( f(u'_i) \) and \( v_i \), we can deduce that \( f(u'_i) = v_i \) for all \( i \). Finally, this yields \( PM^D(u_i, v_i) = k_i \), and taking all the \( f_i = f \) gives all compatible functions, which concludes the proof.

In this proof, we used a constant \( \# \). However, it can also be conducted without using a constant alphabet; indeed, constants can be encoded with parameters, as shown in Appendix B.

Lemma 9 is still valid if \( PM^D \) is replaced by \( FM^D \). This time, we conduct this proof without resorting to the use of constants. This result will be used twice: once for the proof of theorem 17, and again to show that we can once more encode constants into \( \Pi \) using Lemma 25 in Appendix B.

Lemma 10. Let \( u_1, \ldots, u_n \) and \( v_1, \ldots, v_n \) be parameterized words over \( \Pi \) such that \( k_i = |v_i| - |u_i| \geq 0 \), and \( k = \sum_{i=1}^{n} k_i \). Then there exist \( u \) and \( v \) two parameterized words over \( \Pi \cup \{\#\} \), where \( \# \) is a fresh variable, such that the following are equivalent:

1. \( FM^D(u, v) \leq k \)
2. For all \( 1 \leq i \leq n \), \( FM^D(u_i, v_i) \leq k_i \), and the applications \( f_i \) realizing those matchings are all compatible.

Proof. The same technique as Lemma 9 is used but \( u \) and \( v \) are defined as \( u = \#^{k+1} u_1 \# u_2 \# \ldots \# u_n \) and \( v = \#^{k+1} v_1 \# v_2 \# \ldots \# v_n \) where \( \#^{k+1} \) denotes \( k + 1 \) repetitions of the character \( \# \). The full proof can be found in Appendix A.

Finally, the same block result holds for \( FM^D \), and will be used in the proof of theorem 15.

Lemma 11. Let \( u_1, \ldots, u_n \) and \( v_1, \ldots, v_n \) be parameterized words over \( \Pi \) such that for every \( 1 \leq i \leq n \), \( k_i = |u_i| - |v_i| \geq 0 \), and \( k = \sum_{i=1}^{n} k_i \). Then there exist \( u \) and \( v \) two parameterized words over \( \Pi \cup \{\#\} \), where \( \# \) is a fresh variable, such that the following are equivalent:

1. \( FM^D(u, v) \leq k \)
2. For all \( 1 \leq i \leq n \), \( FM^D(u_i, v_i) \leq k_i \), and the applications \( f_i \) realizing those matchings are all compatible.
Proof Idea. The difference with Lemma 10 is that some \( \# \) symbols might be deleted, while some base letters could be mapped to \( \# \). To ensure this does not happen, we define 

\[
    u = \#^N u_1 \#^N u_2 \ldots \#^N u_n \#^N \quad \text{and} \quad v = \#^N v_1 \#^N v_2 \ldots \#^N v_n \#^N.
\]

The full proof can be found in Appendix A.

The technique of block-by-block matching will be used in all the reductions, to encode multiple constraints in a single \( PM \) or \( FM \) instance.

### 4.2 1-to-1 Parameterized Matching \( PM \)

We now focus on the complexity of the \( PM^d \) problems. These problems, as well as function matching problems, are all clearly in \( \text{NP} \): given the list of deletion, insertion or substitution operations to do and the matching to apply, it is easy to check that the solution is correct.

For the reductions in this paper, we always assume that words are written without constants, that is to say \( \Sigma = \emptyset \), since this is sufficient for \( \text{NP} \)-completeness results. This choice is also motivated by the results of Appendix B, which show that \( \Sigma \) can in most cases be coded into \( \Pi \).

\[\iffalse\]

\[\text{Theorem 12. The 1-to-1 Parameterized Matching with deletions } \text{PM}^D \text{ is } \text{NP-complete.}
\]

\[\fi\]

The proof is a reduction from the \( \text{NP} \)-complete problem \( 3\text{-coloring} \)[20]. Given an instance \( G \) of \( 3\text{-coloring} \), we will construct two words \( u \) and \( v \). The word \( v \) will represent the list of vertices and edges of \( G \), while the word \( u \) will list the color of each vertex, and the possible coloring of each pair of vertices joined by an edge. By deleting characters from \( u \), we make a choice for the coloring of each vertex, and thus each edge. The function \( f \) answering the parameterized matching problem will assign a choice of color to each vertex. The instance that we define should be positive iff \( G \) is 3-colorable. More formally:

\[\iffalse\]

\[\text{Proof. The 3-COLOURING problem is defined as follows:}
\]

\[\iffalse\]

\[\text{Input: } G = (V, E) \text{ a graph with vertices } V \text{ and edges } E \]

\[\iffalse\]

\[\text{Output: A coloring } c : V \to \{c_1, c_2, c_3\} \text{ such that for all } \{u, v\} \in E, c(u) \neq c(v) \]

\[\iffalse\]

Let \( G = (V, E) \) be an instance of \( 3\text{-coloring} \), and let \( V = \{x_1, \ldots, x_n\} \) be the set of its \( n \) vertices, and \( E = \{e_1, \ldots, e_m\} \) be the set of its edges. The parameter alphabet \( \Pi \), of polynomial size in \( O(|G|) \) will contain:

- \( x_1, \ldots, x_n \), corresponding to the vertices of \( G \);
- \( n \) copies of the parameters corresponding to the colors \( c_1, c_2, c_3 \); \( c_1, c_2, c_3 \) for \( 1 \leq i \leq n \);
- for every edge \( e \), the delimiters \( Y_e \) and \( \square_e^1, \square_e^2 \);
- \( 2n \) bottom symbols, \( \sqcap_1, \sqcap_2 \) for \( 1 \leq i \leq n \), which will be used to fix the image of some parameters.

First, we define words that will encode the constraint that each vertex is colored, and we make sure that the unused color variables cannot be assigned elsewhere. For \( 1 \leq i \leq n \), \( u_1^i = u_1^i = c_1^i c_2^i c_3^i, u_1^i = x_i \) and \( v_1^i = \sqcap_1^i \sqcap_2^i \). We then define words that include all possible colorings of each edge, and we make sure to use enough brackets. For every edge \( e = \{x_i, x_j\} \), we define \( u_2^e = \square_e^x c_1^i c_2^j c_3^j \square_e^x \square_e^c c_1^i c_2^j c_3^j \square_e^c \square_e^d c_1^j c_2^j c_3^j \square_e^d \square_e^e c_1^i c_2^j c_3^j \square_e^e \) and \( v_2^e = Y_e x_i x_j Y_e \).

Applying Lemma 9 to \( u_1^1, \ldots, u_1^n, u_1^n, u_2^1, \ldots, u_2^m, v_2^1, \ldots, v_2^m \), we obtain \( u \) and \( v \). Let \( k = |u| - |v| = 3n + 20m \). We now show that \( G \) is 3-colorable \( \iff PM_D(u, v) \leq k \).\]

\[\fi\]
Thus Lemma 9 yields \( u \leq 3 \) allowed arrangements of the colors is in \( \text{Theorem 13} \).

We do a simple reduction from \( \text{PM}_D \) as an instance of \( u, v, k \).

**Proof.** We do a simple reduction from \( \text{PM}_D \). Let \( u, v, k \) be an instance of \( \text{PM}_D \). If the instance is trivially false (that is to say, \( k \neq |u| - |v| \)), answer negatively. Else, consider \( u, v, k \) as an instance of \( \text{PM}_D^{\text{DIS}} \). If this is a negative instance for \( \text{PM}_D^{\text{DIS}} \), it is also negative.
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for $PM^D$. Furthermore, if it is a positive instance for $PM^{DIS}$, exactly $k$ deletions should be applied, and so no substitution or insertion are used in that solution. Hence, that solution is also a solution to $PM^D$, and the reduction holds.

The same result in fact holds for all other distances, and in particular the longest common sub-word distance ID. This proves once again the result shown in [26]:

\begin{itemize}
  \item Corollary 14. The problems $PM^I$, $PM^{DI}$, $PM^{IS}$, $PM^{DS}$ are all \textsc{np}-complete.
\end{itemize}

Proof. From Section 3.2, $PM^I$ and $PM^D$ are equivalent in the 1-to-1 case. For the other problem, we do an immediate reduction from $PM^I$ or $PM^D$ analog to the proof of Theorem 13.

We now turn to proofs of \textsc{np}-completeness without the restriction asking $f$ to be injective.

### 4.3 Function Matching $FM^d_1$

The problem considered in this section is the one where both deletions and $f$ are applied to the first word. A reduction very similar to the one given for $PM^D$ is used.

\begin{itemize}
  \item Theorem 15. $FM^d_1$ is \textsc{np}-complete.
\end{itemize}

Proof Idea. The reduction follows the same idea as in Theorem 12. Since the function $f$ realizing the matchings is not injective in this version, it will be used to send every vertex to its color. Moreover, we add more “sink” \bot letters to force the image of every unused letter. The full proof can be found in Appendix A.

This again ensures the \textsc{np}-completeness of the problem for all edit distances, using the same proof as for Theorem 13.

\begin{itemize}
  \item Corollary 16. The problem $FM^{DIS}_1$ of function matching under the Levenshtein distance is \textsc{np}-complete. The problems $FM^I_1$, $FM^{ID}_1$, $FM^{IS}_1$, $FM^{DS}_1$ are all \textsc{np}-complete too.
\end{itemize}

We can notice that the problem $FM^S_1$, where substitution is the only operation allowed, is polynomial-time solvable. Intuitively, for each parameter, consider the possible parameters that it could be mapped to, and their respective number of occurrences. Then, choose the letter with the highest number of occurrences for the mapping. The remaining letters are then substituted.

### 4.4 Function Matching $FM^d_2$

The problem considered in this section is the one where deletions are applied to the second word, while $f$ is applied to the first word.

\begin{itemize}
  \item Theorem 17. $FM^D_2$ is \textsc{np}-complete.
\end{itemize}

Proof Idea. The proof is very similar to the previous case, but the bracketing has to be adapted. Separators $Y^e$ are duplicated enough times to ensure that no vertex variable is mapped to them. The full proof can be found in Appendix A.

\begin{itemize}
  \item Corollary 18. $FM^I_1$, $FM^{DI}_2$, and $FM^L_2$ are all \textsc{np}-complete.
\end{itemize}

Proof. $FM^I_1$ is equivalent to $FM^D_2$. For the two other problems, we use a reduction from $FM^D_2$ exactly like in Corollary 14.
This last result completes the picture of NP-completeness proofs, and indicates that computing the distances between parameterized words defined in Section 3.3 is in general an NP-complete problem.

Similarly to $FM_1^S$, $FM_2^S$ is also polynomial-time solvable.

## 5 Approaches to Solve Parameterized Matching

In this section, we discuss two ways to get around the difficulty of the parameterized matching problems. The first one is to design an FPT algorithm in the alphabet size, and the second one is to translate the problem into a SAT formalism, with the intent of using a SAT-solver.

### 5.1 An FPT Algorithm in the Alphabet Size

The fact that $\Sigma$ and $\Pi$ are part of the input is what makes the various parameterized matching problems NP-hard. When the alphabet size is considered fixed, a simple polynomial algorithm can be used, which generalizes the “naïve” brute force algorithm of Theorem 1 of [26]:

**Algorithm 1** Simple FPT algorithm for $FM^d$.

$m \leftarrow 0$

*for* all functions $f : \Pi \rightarrow \Pi$

$\text{dist} \leftarrow d(f(u), v)$

*if* $\text{dist} \leq m$ then

$m \leftarrow \text{dist}$

*end if*

*end for*

▶ **Theorem 19.** Let $d$ be a distance. Algorithm 1 computes $FM^d(u, v)$ in time $O(|\Pi||\Pi||u||v|)$

**Proof.** Algorithm 1 uses an exhaustive search and finds $\min_{f : \Pi \rightarrow \Pi} d(f(u), v)$, which is the definition of $FM^d(u, v)$. Furthermore, there are $|\Pi|^2$ functions from $\Pi$ to $\Pi$, and computing $d(f(u), v)$ is done in time $O(|f(u)||v|) = O(|u||v|)$, hence a total running time in $O(|\Pi|^2|u||v|)$. ◀

Note that this also leads to a similar algorithm for $PM^d$ by iterating over injective functions rather than all functions from $\Pi$ to $\Pi$.

### 5.2 A MaxSat Formulation of Parameterized Matching

In this section, we encode $PM^d$ problems into SAT problems, with the intent of solving them with a SAT solver. More precisely, we will use the weighted max-SAT variant of SAT, which is defined in the following way:

- **Input:** a set $V$ of literals, a formula $\varphi = \bigwedge_{i=1}^{n} \varphi_i$ on $V$ in conjunctive normal form (CNF),
  a weight function $w : [1, n] \rightarrow \mathbb{N}$.
- **Output:** a valuation $\nu : V \rightarrow \{0, 1\}$ such that $\sum_{\nu \models \varphi_i} w(i)$ is maximal.

Moreover, we will sometimes use a partially weighted variant of Max-SAT, which is defined in the following way:
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- **Input:** a set $V$ of literals, a satisfiable formula $\varphi_c$ on $V$ in CNF, a formula $\varphi_w = \bigwedge_{i=1}^n \varphi_i$ on $V$ in CNF and a weight function $w : [1,n] \to \mathbb{N}$.
- **Output:** a valuation $\nu : V \to \{0,1\}$ such that $\nu \models \varphi_c$ and $\sum_{\nu \models \varphi_i} w(i)$ is maximal.

In that case, clauses of $\varphi_c$ are called “hard” clauses while clauses of $\varphi_w$ are called “soft clauses”. We give a proof of the equivalence in Proposition 26 of Appendix C.

We will define an encoding of an instance $(u, v)$ of $PM^d$ such that an assignment of the variables of $V$ will define an alignment between $u$ and $v$. First, we make a link between the ID edit distance and the length of the optimal alignment between two strings.

**Definition 20.** Let $u$ and $v$ be two words on $\Pi$, such that $p = |u|$ and $p' = |v|$. A set $A \subset [1,|u|] \times [1,|v|]$ is an alignment between $u$ and $v$ iff the following are true:

1. Each position of $u$ appears at most once: For all $1 \leq i \leq p$ and $1 \leq j, j' \leq p'$, if $(i, j) \in A$ and $(i, j') \in A$, then $j = j'$.
2. Each position of $v$ appears at most once: For all $1 \leq j \leq p'$ and $1 \leq i, i' \leq p$, if $(i, j) \in A$ and $(i', j) \in A$, then $i = i'$.
3. There are no crossings: if $(i, j) \in A, (i', j') \in A$, and $i' > i$, then $j' > j$.
4. Aligned positions match in $u$ and $v$: if $(i, j) \in A$, then $u_i = v_j$.

An alignment relates to the insertion/deletion distance $ID$ in the following way:

**Theorem 21.** Let $u, v$ be words on $\Pi$ and $k \leq |u| + |v|$ be an integer. The following are equivalent:

1. There exists an alignment $A$ such that $2|A| = |u| + |v| - k$.
2. $ID(u, v) \leq k$.

**Proof.** The proof, which works by induction, can be found in Appendix C.

We now turn to the max-SAT encoding of our problem.

**Theorem 22.** Let $u$ and $v$ be two words over $\Pi$. There exists a formula $\varphi_{u,v} = \varphi_c \land \varphi_w$ and a weight function $w$, instance of the partially weighted Max-SAT problem such that the following are equivalent:

- $\nu$ is a solution to this partially weighted Max-SAT instance and satisfies $k$ clauses of $\varphi_w$.
- There exists a function $f : \Pi \to \Pi$ and an alignment between $f(u)$ and $v$ of size $k$.

The formula $\varphi$ uses $|\Pi||p| + |\Pi|^2$ variables and is of size $O(m^2p^2)$, where $m = |u|$ and $p = |v|$. Moreover, there exists $\varphi^{inj}$ of size $O(|\Pi|^3)$ such that the above result is true for $f$ injective by replacing $\varphi_c$ with $\varphi'_c = \varphi_c \land \varphi^{inj}$.

In particular, finding the valuation maximizing $k$ gives a maximal alignment between $u$ and $v$, and with Theorem 21, the distance $ID(u, v)$.

**Proof.** For this proof, we fix an ordering on the alphabet $\Pi = \{a_1, \ldots, a_n\}$. We define the set of literals $V$ as $V = \{x_{i,j} \mid 1 \leq i \leq |u|, 1 \leq j \leq |v|\} \cup \{y_{a,b} \mid a \in \Pi, b \in \Pi\}$. Intuitively, $x_{i,j}$ represents a match between position $i$ and $j$ in the alignment, and $y_{a,b}$ will represent the fact that $f(a) = b$. We define the following sets of formulas, where all indices $i$ are taken between 1 and $m$ and all $j$ between 1 and $p$, and $a$ and $b$ are taken in $\Pi$:
We then define \( \varphi_c \) as the conjunction of all the formulas (NoDouble i), (NoDouble j), (NoCrossing), (Function), and (Match). Furthermore, we define \( \varphi^{inj} \) as the conjunction of all the (Injectivity) formulas. Lastly, we define \( \varphi_w = \bigwedge_{1 \leq i \leq m} \varphi^3_i \), and set \( w(C) = 1 \) for every clause \( C \) of \( \varphi_w \).

There are \( m \binom{2}{0} \) (NoDouble i) formulas, \( p \binom{m}{n} \) (NoDouble j), \( \binom{m}{n} \binom{m}{n} \) (NoCrossing), \( n \binom{2}{0} \) (Function) and (Injectivity) formulas, \( pm \) (Match) formulas and \( n \) (ExistsMatch) formulas.

We now prove both implications of the theorem. Suppose \( \nu \) is a valuation satisfying \( \varphi_c \) and \( k \) clauses of \( \varphi_w \). We define, for all \( a, b \in P_i \), \( f(a) = b \) if and only if \( \nu(y_{a,b}) = \top \). Since \( \nu \) satisfies all the (Function) formulas, this is a correct definition of a (partial) function. We define \( A = \{(i, j) \mid \nu(x_{i,j}) = \top\} \). \( A \) is an alignment between \( f(u) \) and \( v \). Indeed: (NoDouble i) and (NoDouble j) ensures point 1 and 2. of Definition 20, (NoCrossing) ensures point 3., and Match ensures point 4. The size of \( A \) is the number of \( x_{i,j} \) instantiated to \( \top \), which is exactly the number of clauses of \( \varphi_c \), satisfied, i.e., \( k \).

Suppose now that there exists a function \( \Pi \rightarrow \Pi \) and an alignment \( A \) between \( f(u) \) and \( v \). Similarly, we define \( \nu(y_{a,b}) = \top \) if and only if \( f(a) = b \), and \( \nu(x_{i,j}) = \top \) if and only if \( (i,j) \in A \). Since \( A \) is an alignment, \( \nu \) satisfies (NoDouble i), (NoDouble j), and (NoCrossing). Since \( f \) is a function, (Function) is satisfied. Finally, if \( \nu(x_{i,j}) = \top \), then \( (i,j) \in A \), and since \( A \) is a matching, \( f(u)_i = f(u)_i = v_j \) and \( \nu(y_{u_i,v_j}) = \top \).

The proof for \( \varphi^{b} \) is the same, and (Injectivity) ensures the injectivity of \( f \). □

What is more, this proof can be adapted to change the ID distance to the Levenshtein distance, simply by choosing to consider all the (Match) formulas as soft clauses.

6 Experiments

The two approaches presented in Section 5 were implemented in Python to solve PMID. They are available under the GPL license at https://github.com/AaronFive/paramatch. The FPT algorithm of Section 5.1 is implemented in the function parameterizedAlignment of file fpt_alphabet_size.py. The MaxSAT-reduction of Section 5.2 is implemented in the function make_sat_instance of file sat_instance.py. The MaxHS solver [18] available at http://www.maxhs.org is used by our script to solve the MaxSAT instances derived from the PMID instances.

Our initial motivation to introduce parameterized matching under various distances is theater play comparison. To represent the structure of a theater play, we represent each character by a letter of the alphabet, and create the parameterized word obtained by considering the succession of all consecutive speakers. To check their adequacy with real data,
we use a corpus of theater plays in which each character is represented by one letter of the alphabet, and each act of the play is represented by a string corresponding to the sequence of speaking characters. A letter may be duplicated in this string if the corresponding characters has lines in the end of a scene and in the beginning of the next one. Therefore, the edit distance between two parameter words representing acts will be small if both acts have a similar structure in terms of succession of speaking characters. We selected a corpus of 10 pairs of plays where one inspired the other, and performed 47 comparisons between pairs of acts. Among those comparisons, 26 were solved by the maxSAT algorithm and all by the FPT algorithm (detailed results are presented in the supplementary material available at https://github.com/AaronFive/paramatch/tree/main/corpus10pairs), with a 800 second timeout. The computation times are obtained on a XMG laptop running on Windows, with a 2.60 Ghz processor and 16 Gb RAM. Only the running time of MaxHS is provided, the encoding into a MaxSAT formula usually runs in approximately 1 second. Note that all instances are solved faster by the FPT algorithm than by the MaxSAT approach. The analysis of running times depending on the product of the lengths of the input strings (see supplementary material) shows that the MaxSAT approach may be relevant for strings with more than 10 distinct characters, but where the product of the length of input strings may not exceed 2000.

7 Conclusion

In this paper, we studied the complexity of several variants of the edit distance problem between parameterized words. We proved the NP-completeness of all previously unsolved cases, including the Levenshtein distance left open in [24], and provided practical approaches to solve real instances of those problems. We also studied similar problems for various definitions of words with parameters, namely parameter words and parameterized expressions, proving some relationships with parameterized word problems.

As future work, we will study the restrictions introduced in [21, 22] for a pattern matching problem with patterns in the parameter, in order to obtain polynomial time algorithms for the edit distance between parameterized words. Moreover, we will explore the question of distance between sets of words, in particular when they are defined through generalizations of automata. These problems are variants of the notion of distance between regular languages as defined in [12]. In this context, we can notice that different notions of automata can be considered: either automata generating parameterized words, or automata using parameters to define languages over classical words, with two different semantics as defined in [11].

References


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Details of the proofs

Proof of Lemma 7. We proceed by induction on \( k \). If \( k = 0 \), then \( u \) and \( v \) are parameterized matching, and \( f(u) = v \), thus \( d(f(u), v) = 0 \). Suppose the result holds until a fixed \( k \). Suppose \( PM^d(u, v) = k + 1 \). There exist \( f, u'' \) and \( u' \) such that \( d(u, u'') = 1 \), \( d(u'', u') = k \), and \( f(u') = v \). Hence \( PM^d(u'', v) \leq k \), and by induction hypothesis \( d(f(u''), v) \leq k \). Moreover,
since \( d(u, u'') = 1 \), we get \( u'' \) from \( u \) by applying only one operation. We prove that regardless of this operation, \( d(f(u), f(u'')) = 1 \), and thus \( d(f(u), v) \leq d(f(u), f(u'')) + d(f(u''), v) \leq k+1 \) which will conclude the proof. There are 3 cases to consider:

- If the operation is a deletion, \( u = v_1xv_2 \) and \( u'' = v_1v_2 \) for some words \( v_1 \) and \( v_2 \) and some letter \( x \). Then \( f(u) = f(v_1)f(x)f(v_2) \) and we can obtain \( f(v_1)f(v_2) = f(u'') \) by deleting \( f(x) \).
- If it is an insertion, \( u = v_1v_2 \) and \( u'' = v_1xv_2 \), and we can similarly go from \( f(u) \) to \( f(u'') \) by inserting \( f(x) \).
- If it is a substitution, \( u = v_1xv_2 \) and \( u'' = v_1yv_2 \), and we can go from \( f(u) \) to \( f(u'') \) by replacing \( f(x) \) with \( f(y) \).

Hence \( d(f(u), f(u'')) = 1 \), which concludes the proof for \( PM^d \).

Since this proof does not use the fact that \( f \) is 1-to-1, it also stands for \( FM^d_1 \).

**Proof of Lemma 8.** It is obvious that \( d(zu, zv) \leq d(u, v) \), so we only prove \( d(u, v) \leq d(zu, zv) \). We prove that any rewriting sequence from \( zu \) to \( zv \) can be modified such that no edit operation is applied in \( z \). This will be enough to prove the result, as the edit sequence obtained can be seen as an edit sequence between \( u \) and \( v \). We proceed by induction on the size of \( z \). Suppose \( |z| = 1 \). Then \( z = a \in \Sigma \cup \Pi \). We can consider that no character is modified twice in an edit sequence (i.e. no character is inserted and then deleted, or inserted and then substituted etc.), as that is always sub-optimal. Suppose \( z \) is modified. There are 3 possible cases:

1. There is an insertion in \( z \), hence a word \( w \) ends up being inserted before \( a \). Since \( zv = av \) starts with \( a \), \( w \) must start with an \( a \), hence \( w = aw' \). We insert \( w' \) to the right of \( z \) instead with the same operations. If \( z \) should be deleted or substituted, we apply the same operation to the new \( a \) instead. These operations yield the same result, and do not modify \( z \).

2. There is a deletion in \( z \), and hence \( a \) is deleted. Since this an optimal rewriting sequence, no \( a \) is created at that position through insertion or substitution afterwards. Since \( av \) starts with an \( a \), \( u \) must be of the form \( u = sau' \), where all the characters in \( s \) are deleted, and \( a \) isn’t. Deleting \( sa \) instead of \( as \) yields the same result, and doesn’t modify \( z \).

3. There is a substitution in \( z \), hence \( a \) is modified into a character \( b \neq a \), that will not be further modified. Since \( av \) starts with \( a \), \( an \) has to be inserted in \( z \), which is handled in case 1.

Hence, we can consider that every edit operation is done in \( u \), and \( d(au, av) = d(u, v) \). Suppose now that the result is proven for \( |z| = k \), and let \( z = az' \), with \( |z'| = k \). Using the base case and the case for \( |z| = k \), we have \( d(zu, zv) = d(azu, azv) = d(zu, zv) = d(u, v) \), which concludes the proof.

**Proof of Lemma 10.** Let \( # \) be a fresh parameterized letter. Let then \( u = #^{k+1}u_1#u_2#\ldots #u_n \) and \( v = #^{k+1}v_1#v_2#\ldots #v_n \), where \( #^{k+1} \) denotes \( k+1 \) repetitions of the character \( # \). The proof of the reverse direction is the same as in Lemma 9, so we only prove the other one.

Assume \( FM^P_m(u, v) \leq k \). Let \( v' \) and \( f \) realize this parameterized match.

We prove that \( f(#) = # \), and that no other character is sent to \( # \) by \( f \). Indeed, \( v \) starts with \( k+1 \) symbols \( # \), which ensure that \( v' \) starts with the letter \( # \). Since \( u \) starts with \( # \) and \( f(u) = v' \), \( f(#) = # \). Furthermore, this implies that since \( |u|_# = k+n \), \( |f(u)|_# = |v'|_# \geq k+n \). Since \( v' \) is obtained from \( v \) by deletions, we have \( |v'|_# \leq |v|_# = k+n \). Hence \( |v|_# = k+n \) and all those inequalities are equalities, which is only the case when no \( # \) symbols is deleted from \( v \), and that for all \( x \neq # \), \( f(x) \neq # \).
Since all the # symbols are left untouched, the rest of the proof is the same as in Lemma 9, and all of the factors \( u_i \) and \( v_i \) are parameterized matching.

**Proof of Lemma 11.** Let \( \# \) be a fresh parameterized letter, and \( N = k + 2 \).

Let then \( u = \#^N u_1 \#^N u_2 \ldots \#^N u_n \#^N \) and \( v = \#^N v_1 \#^N v_2 \ldots \#^N v_n \#^N \). Once again, we only prove the non-trivial implication.

Suppose \( \text{FM}_1^D(u,v) \leq k \), and let \( f \) and \( u' \) realize this matching. Since \( u \) starts with \( k + 1 \) copies of \( \# \), \( u' \) starts with \( \# \). Since \( v \) starts with \( \# \) too, \( f(\#) = \# \).

We now prove that we can consider that for all \( x \neq \# \), \( f(x) \neq \# \). This will also imply that no \( \# \) symbol is deleted from \( u \). Let \( S = \{ a \in \Pi \mid f(a) = \# \} \) be the set of symbols (different from \( \# \)) sent to \( \# \). Since \( |u|_\# = |v|_\# \), the number of deleted \( \# \) symbols from \( u \) is exactly \( |u|_S \), hence \( |u|_S \leq k \). Let us now consider the leftmost occurrence of an element of \( S \) in \( u' \), that we denote by \( a \). The letter \( a \) appears in \( u \) in a factor of the form \( \#^N w_1 a w_2 \#^N \). Since all \( \# \) in \( v \) appear in blocks of size \( N \), \( a \) must contribute to such a block, after deletions and application of \( f \). We distinguish two cases:

1. The entirety of the word \( w_1 \) is deleted. In this case, at least one symbol \( \# \) from the left \( \#^N \) block is deleted; otherwise \( f(\#^N) f(a) = \#^{N+1} \) would be a factor of \( v \), which is impossible. Thus, choosing not to delete \( \# \) and to delete \( a \) instead yields the same result.
2. \( w_1 \) is not deleted. Since no character from \( S \) appears to the left of \( a \), \( f(a) \) is the start of a \( \#^N \) block. Furthermore, since \( |u|_S \leq k \), it is not possible to form \( \#^N \) with only \( a \) and \( w_2 \), and characters from the right \( \#^N \) contribute to it. Hence, at least one \( \# \) symbol from this right block is deleted. Like before, the same result can be obtained by not deleting it, and deleting \( a \) instead.

Either way, we can repeat this process to eliminate all occurrences of characters of \( S \) and of deletions of \( \# \), which proves that we can consider that for all \( x \neq \# \), \( f(x) \neq \# \). Once again, we are taken back to the conditions of Lemma 9, and the rest of the proof follows.

**Proof of Theorem 15.** We define \( \Pi \) like in Theorem 12, and we add the letters \( \bot_1, \bot_2, \bot_3, \bot_4 \) and \( \bot_5 \). Similarly, we define \( u_1^\epsilon, v_1^\epsilon, u_1^\bot, v_1^\bot, u_2^\epsilon, v_2^\epsilon, u_2^\bot, v_2^\bot, \) and \( v_2^\bot \) just like in Theorem 12. Additionally, we define for every edge \( e \),

\[
\begin{align*}
  u_1^\epsilon &= \bot_1 \bot_2 \bot_3 \bot_4 \bot_5 \text{ and } v_1^\epsilon &= \bot_1 \bot_2 \bot_3 \bot_4 \bot_5. \\
  \text{We then apply Lemma 11 with} \\
  u_1^1, u_1^2, \ldots, u_1^n, u_2^1, u_2^2, \ldots, u_2^m, u_1^1, \ldots, u_1^m \\
  \text{and} \\
  u_1^1, u_1^2, \ldots, u_1^n, v_2^1, v_2^2, \ldots, v_2^m, v_1^1, \ldots, v_1^m \\
\end{align*}
\]

to obtain \( u, v, \) and \( k \). We show that \( G \) is 3-colorable \( \Rightarrow \text{FM}_1^D(u,v) \leq k \).

\( \Rightarrow \) Suppose \( G \) is 3 colorable. Define \( f \) like in Theorem 12 on the \( c_i^e \) and \( \Box_y^e \). Let \( e \) be an edge and \( k_e \in [1,6] \) be the integer such that \( f(\Box_y^e) \) is defined. We map every remaining \( \Box_y^e \) in the following way:

\[
f(\Box_y^e) = \begin{cases} 
\bot_1 & \text{if } i < k_e, \\
Y^e & \text{if } i = k_e, \\
\bot_{i-1} & \text{if } i > k_e.
\end{cases}
\]  \hspace{1cm} (1)

It is then easy to check that \( d(f(u),v) = k \), and thus \( \text{FM}_1^D(u,v) \leq k \).
We define for $1 \leq i \leq n$ and $1 \leq t \leq 3$, $\text{col}(c_t^i) = c_i$. If $x_i$ is a vertex of $G$, define $c(x_i)$ to be $\text{col}(c_{x_i}^i)$, where $c_{x_i}^i$ is the only element such that $f(c_{x_i}^i) = x_i$. We show in what follows that (1) this function definition is correct and (2) it is a valid coloring, i.e. if $e = \{x_i, x_j\}$ is an edge, $c(x_i) \neq c(x_j)$.

(1): The same points 1. and 2. from the proof of Theorem 12 apply, hence for every $1 \leq i \leq n$, exactly one element from $\{c_1^i, c_2^i, c_3^i\}$ is sent to $x_i$, while the two others are sent to $\perp_1$ and $\perp_2$, hence the result.

(2): Let $e$ be an edge. The words $u_1^e$ and $v_1^e$ are in matching, which is done with exactly one deletion. Hence, there exists $k_e$ such that

$$f(\square_t^i) = \begin{cases} \perp_i & \text{if } i < k_e, \\ \perp_{i-1} & \text{if } i > k_e. \end{cases}$$

(2)

Moreover, $u_2^e$ and $v_2^e$ are in matching. Since $Y_e$ appears in $v_2^e$ and all the characters in $u_2^e$ apart from $\square_{k_e}^i$ have an image different from $Y_e$, $f(\square_{k_e}^i) = Y_e$. Hence, the only characters that are not suppressed from $u_2^e$ are the two characters between the $\square_{k_e}^i$. Denoting them by $c$ and $c'$, the construction of the word ensures that $\text{col}(c) \neq \text{col}(c')$. Hence, if $e = \{x_i, x_j\}$, we have proven $c(x_i) \neq c(x_j)$, which is (2).

The coloring $c$ is therefore valid, which concludes the proof.

\textbf{Proof of Theorem 17.} Let $G = (V, E)$, with $V = \{x_1, \ldots, x_n\}$ and $\{e_1, \ldots, e_m\}$. Like in the 1-to-1 case, we construct factors $u_i$ and $v_i$ to encode vertex coloring. The parameter alphabet contains:

- $x_1, \ldots, x_n$, corresponding to $V$,
- the colors $c_1, c_2, c_3$,
- for every $e \in E$, the delimiters $Y_e$,
- for every $e \in E$ and every $1 \leq i, j \leq 3$, $i \neq j$, the delimiters $Y_{ij}^e$.

We define for $1 \leq i \leq n$, $u_1^i = x_i$ and $v_1^i = c_1 c_2 c_3$. If $e$ is an edge and $c_i$ and $c_j$ are two colors, we denote $w^e(c_i, c_j) = Y_{ij}^e Y_e Y_e Y_{ij}^e$. For every edge $e = \{x_i, x_j\}$, we now define $u_2^e = Y_e Y_e Y_e Y_e x_i Y_e Y_e Y_e Y_e$ and $v_2^e = w^e(c_1, c_2) w^e(c_2, c_1) w^e(c_2, c_3) w^e(c_3, c_1) w^e(c_3, c_2)$.

We now apply Lemma 10 with $u_1^i = u_2^1 \ldots u_2^i \ldots u_2^n$, $v_1^i = v_2^1 \ldots v_2^i \ldots v_2^n$, to obtain $u$ and $v$.

$\Rightarrow$ Suppose $G$ is 3-colorable, and let $c : V \rightarrow \{c_1, c_2, c_3\}$ be a valid coloring. Define $f \upharpoonright V = c$. For every edge $e = \{x_i, x_j\}$, let $s$ and $t$ be such that $c(x_i) = c_s$ and $c(x_j) = c_t$. We then define $f(Y_e) = Y_{s+t}$. It is easy to check now that $d(f(u), v) = k$, and hence $FM_P^P(u, v) \leq k$.

$\Leftarrow$ Suppose now that $FM_P^P(u, v) \leq k$. We will show that $f \upharpoonright V$ defines a 3-coloring of $G$, by showing that (1) for all $x \in V$, $f(x) \in \{c_1, c_2, c_3\}$ and (2) If $\{x, y\} \in E$, then $f(x) \neq f(y)$.

- Lemma 10 ensures that the words $u_i$ and $v_i$ are in matching, which proves (1).
- Lemma 10 also ensures that the words $w^e$ and $\varepsilon^e$ are in matching. Let $e \in E$, with $e = x_s, x_t$. We have $|u_2^e| = 6$, hence $|f(u_2^e)| = 6$. Since $c_1, c_2$ and $c_3$ each occur exactly 4 times in $v_2^e$, they cannot occur 6 times after deletions, and $f(Y_e) \notin \{c_1, c_2, c_3\}$. Hence, there exist $i \neq j$ with $1 \leq i, j \leq 3$ such that $f(Y_e) = Y_{ij}^e$. This implies that all but one of the $w^e$ factors from $v_2^e$ are suppressed, and that the remaining one is $w^e(c_i, c_j)$. Hence $f(x_s) = c_i$ and $f(x_t) = c_j$, which proves (2).
B Encoding Constant Alphabet $\Sigma$ in $\Pi$

We show why it is always possible to consider that $\Sigma = \emptyset$ for certain problems. These results use the lemmas proved in Section 4.1.

**Lemma 23.** Let $d$ be a distance, $k$ an integer and $u$ and $v$ be two parameterized words over the alphabet of constants $\Sigma$ and the alphabet of parameters $\Pi$. There exist words $\tilde{u}$ and $\tilde{v}$ over the alphabet of constants $\emptyset$ and the alphabet of parameters $\Pi' = \Pi \cup \Sigma$ such that the following are equivalent:

- $PM^d(u, v, k)$ is realized by $f$;
- $PM^d(\tilde{u}, \tilde{v}, k)$ is realized by $f$.

In particular, this implies that if $PM^d(\tilde{u}, \tilde{v}) \leq k$, all functions $f$ realizing this matching verify that for all $x \in \Sigma$, $f(x) = x$, and for all $x \in \Pi, f(x) \in \Pi$.

**Proof.** Let $N = k + 1$. If $\Sigma = \{a_1, \ldots, a_n\}$, we define $z$ to be $a_1^N a_2^N \ldots a_n^N u$ and $\tilde{u} = zu$, $\tilde{v} = zv$. It is clear that if $PM^d(u, v) \leq k$ then $PM^d(\tilde{u}, \tilde{v}) \leq k$, by following the same operations, and applying the same renaming function.

Suppose now that $PM^d(\tilde{u}, \tilde{v}) \leq k$, and let $f$ and $u'$ realize it. Let $i \in [1, n]$. All the letters of $u$ between position $Ni$ and $N(i + 1)$ are $a_i$. At most $k$ of these positions can be modified with an edit operation. Since $N > k$, at least one of these positions is not modified, and thus there exists $j \in [Ni, N(i + 1)]$ such that $u'_j = a_i$. Since all letters in $v$ between position $Ni$ and $N(i + 1)$ are $a_i$, in particular $v_j = a_i$, and hence $f(a_i) = a_i$. This proves that for all $x \in \Sigma$, $f(x) = x$, and thus $f(z) = z$. Since $f$ is 1-to-1, this entails $f(\Pi) \subseteq \Pi$. By Lemma 7, $d(f(\tilde{u}), \tilde{v}) \leq k$. Hence $d(f(zu), zv) = d(zf(u), zv) \leq k$ and by Lemma 8, $d(f(u), v) \leq k$. Hence $PM^d(u, v) \leq k$.

**Remark 24.** Note that the words $\tilde{u}$ and $\tilde{v}$ have a size increased by $N\Sigma$. If less operations are considered, it is possible to reduce this overhead. For example, in the case of $PM^D$, we can take $z$ to be of the form $a_1 \ldots a_n z^N$, to reduce the overhead to $N + \Sigma$.

Similarly, constants can be encoded in $\Pi$ in some $FM$ problems. We prove this result for $FM^2_D$, with the help of the block decomposition allowed by Lemma 10.

**Lemma 25.** Let $u$ and $v$ be two parameterized words over the alphabet of constants $\Sigma$ and the alphabet of parameters $\Pi$. There exist words $\tilde{u}$ and $\tilde{v}$ over the alphabet of constants $\emptyset$ and the alphabet of parameters $\Pi' = \Pi \cup \Sigma$ such that the following are equivalent:

- $FM^2_D(u, v, |v| - |u|)$ is realized by $f$;
- $FM^2_D(\tilde{u}, \tilde{v}, |\tilde{v}| - |\tilde{u}|)$ is realized by $f$.

**Proof.** We write $\Sigma = \{a_1, \ldots, a_n\}$ and $\Pi = \{b_1, \ldots, b_m\}$. We define $z_\Sigma = a_1 \ldots a_n$, and $z_\Pi = b_1 \ldots b_m$. Let $\tilde{u}$ and $\tilde{v}$ be the words obtained by applying Lemma 10 to $z_\Sigma, b_1, b_2, \ldots, b_m, u$ and $z_\Sigma, z_\Pi, z_\Pi, \ldots, z_\Pi, v$. If $FM^2_D(u, v, k)$ is realized by a function $f$, it realizes $FM^2_D(\tilde{u}, \tilde{v}, |\tilde{v}| - |\tilde{u}|)$ too. Indeed, it is enough to apply the same operations in $v$, and to delete all the characters but $f(b_i)$ in the $i$-th copy of $z_\Pi$.

Suppose now that $FM^2_D(\tilde{u}, \tilde{v}) \leq k$, and let $f$ realize it. Then, by Lemma 10, we have:

- $D(z, f(z)) = 0$, and hence $f(z) = z$, which implies that for all $x \in \Sigma$, $f(x) = x$.
- For every $1 \leq i \leq m$, $D(z_\Pi, f(b_i)) = |\Pi| - 1$. Hence $f(b_i)$ is a character of $z_\Pi$, which is some character $b_j \in \Pi$.
- $D(v, f(u)) \leq k$.

Hence $f$ verifies $D(f(v), u) \leq k$ and respects the conditions on $\Pi$ and $\Sigma$, which implies that is also realizes $FM^2_D(u, v, k)$. ▶
The overhead to pay for this transformation is $O(|\Sigma| + |\Pi|^2 + k)$, where the term in $k$ comes from the proof of Lemma 10.

Transposing the technique used for Lemma 25 is not sufficient to get a similar result for $FM_1^{\Pi}$. The question thus remains open in this context.

## C Proofs Regarding the Max-SAT Encoding

**Proof of theorem 21.** We proceed by induction on $|u| + |v|$. If $|u| + |v| = 0$, both $u$ and $v$ are the empty string, and the equivalence is trivial. Fix $n \in \mathcal{N}$ and suppose now that the result holds up for all words $u, v$ such that $|u| + |v| \leq n - 1$. Let $u$ and $v$ be two words such that $|u| + |v| \leq n$. Without loss of generality, consider $|u| \geq |v|$.

Suppose $ID(u, v) \leq k$. Let $\rho$ be a rewriting sequence between $u$ and $v$ of length $k$. If there is no deletion in $u$ in $\rho$, there are only insertions in $v$, and $v$ is a sub-word of $u$, and there exists another rewriting sequence $\rho'$ only deleting letters from $u$. Hence, we can consider that there is at least a deletion in $u$ in $\rho$. Let $p$ be a position at which such a deletion occur, and let $a = u_p$. The word $u$ can be written as $u = u'u''$ for some words $u'$ and $u''$. Define $w = u'u''$. It holds that $d(w, v) \leq k - 1$ and $|w| = |u| - 1$. By induction, there exists an alignment $A$ between $w$ and $v$ such that $2|A| = |w| + |v| - (k - 1) = |u| + |v| - k$. We define $r(i) = \begin{cases} i & \text{if } i < p \\ i - 1 & \text{if } i \geq p \end{cases}$, and $B = \{(r(i), j) \mid (i, j) \in A\}$. Since $A$ is an alignment, so is $B$: it satisfies conditions 1 to 3 of Definition 20, and since $w_r(i) = u_i$, it also satisfies condition 4. Finally, $|B| = |A|$, hence $2|B| = |u| + |v| - k$, hence the result.

Suppose now that there exists an alignment $A$ such that $2|A| = |u| + |v| - k$. Similarly, consider $p$, a position in $u$ such that there does not exist a $j$ with $(p, j) \in A$. If no such position exist, since $|u| \geq |v|$, $u = v$ and the result is proven. Consider $w$ the word obtained by deleting $u_p$ from $u$. It then holds that $|w| = |u| - 1$ and that $2|A| = |u| + |v| - k = |w| + |v| - (k - 1)$.

Defining $B$ in the same way as above yields an alignment between $w$ and $v$ of the same size, and thus by induction, $d(w, v) \leq k - 1$, and since $d(u, w) = 1$, $d(u, v) \leq k$.

**Proposition 26.** Weighted Max-SAT and partial weighted Max-SAT are equivalent.

**Proof.** Encoding a weighted Max-SAT instance as a partially weighted Max-SAT instance is straightforward, as we just have to choose $\varphi_c$ to be empty.

Conversely, given a satisfiable CNF formula $\varphi_c$, a CNF formula $\varphi_w$, and a weight function $w$ on the clauses of $\varphi_w$, we can define a weighted Max-Sat instance in the following way:

- We define $\varphi = \varphi_c \land \varphi_w$.
- We set $W = 1 + \sum_{C \in \text{clause of } \varphi_c} w(C)$, and extend $w$ to clauses of $\varphi_c$ such that $w(C) = W$ for all clauses $C$ of $\varphi_c$.

If $\nu$ is a valuation, we denote by $w(\nu)$ the sum of the weights of all clauses it satisfies $\sum_{C \in \text{clause of } \varphi_c} w(C)$. Since $\varphi_c$ is satisfiable, there exists a valuation $\nu_c$ such that $\nu_c \models \varphi_c$, and $w(\nu_c) \geq |\varphi_c|W$. Let now $\nu$ be a valuation no satisfying a clause of $\varphi_c$. Then $w(\nu) \leq (|\varphi_c| - 1)W + (W - 1) < w(\nu_c)$, hence $nu_c$ is not maximal and cannot be a solution to the weighted Max-SAT instance.
Parameterized Algorithms for String Matching to DAGs: Funnels and Beyond

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Abstract

The problem of String Matching to Labeled Graphs (SMLG) asks to find all the paths in a labeled graph $G = (V, E)$ whose spellings match that of an input string $S \in \Sigma^m$. SMLG can be solved in quadratic $O(m|E|)$ time [Amir et al., JALG 2000], which was proven to be optimal by a recent lower bound conditioned on SETH [Equi et al., ICALP 2019]. The lower bound states that no strongly subquadratic time algorithm exists, even if restricted to directed acyclic graphs (DAGs).

In this work we present the first parameterized algorithms for SMLG on DAGs. Our parameters capture the topological structure of $G$. All our results are derived from a generalization of the Knuth-Morris-Pratt algorithm [Park and Kim, CPM 1995] optimized to work in time proportional to the number of prefix-incomparable matches.

To obtain the parameterization in the topological structure of $G$, we first study a special class of DAGs called funnels [Millani et al., JCO 2020] and generalize them to $k$-funnels and the class $ST_k$. We present several novel characterizations and algorithmic contributions on both funnels and their generalizations.

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Keywords and phrases string matching, parameterized algorithms, FPT inside P, string algorithms, graph algorithms, directed acyclic graphs, labeled graphs, funnels

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1 Introduction

Given a labeled graph $G = (V, E)$ and a string $S$ of length $m$ over an alphabet $\Sigma$ of size $\sigma$, the problem of String Matching to Labeled Graph (SMLG) asks to find all paths in $G$ spelling $S$ in their characters; such paths are known as occurrences or matches of $S$ in $G$. This problem is a generalization of the classical string matching (SM) to a text $T$ of length $n$, which can be encoded as an SMLG instance with a path labeled with $T$. Labeled graphs are present in many areas such as information retrieval [28, 73, 14], graph databases [11, 10, 75, 16] and bioinformatics [27], and SMLG is a primitive operation to locate information on them.

It is a textbook result [2, 29, 81] that the classical SM can be solved in linear $O(n + m)$ time. For example, the well-known Knuth-Morris-Pratt algorithm (KMP) [60] preprocesses $S$ and then scans $T$ while maintaining the longest matching prefix of $S$. However, for SMLG a recent result [12, 34] shows that there is no strongly subquadratic $O(m^{1-\epsilon}|E|), O(m|E|^{1-\epsilon})$ time algorithm unless the strong exponential time hypothesis (SETH) fails, and the most

1 We consider the case where each vertex is labeled with a single character from $\Sigma$. © Manuel Cáceres; licensed under Creative Commons License CC-BY 4.0

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efficient current solutions \[8, 72, 77, 52\] match this bound, thus being optimal in this sense. Moreover, these algorithms solve the approximate version of SMLG (errors in \(S\) only) showing that both problems are equally hard under SETH, which is not the case for SM \[13\].

The history of (exact) SMLG. SMLG can be traced back to the publications of Manber and Wu \[67\] and Dubiner et al. \[33\] where the problem is defined for the first time, and solved in linear time on directed trees by using an extension of KMP. Later Akutsu \[4\] used a sampling on \(V\) and a suffix tree of \(S\) to solve the problem on (undirected) trees in linear time and Park and Kim \[74\] obtained an \(O(N + m|E|)^2\) time algorithm for directed acyclic graphs (DAGs) by extending KMP on a topological ordering of \(G\) (we call this the DAG algorithm). Finally, Amir et al. \[8\] showed an algorithm with the same running time for general graphs with a simple and elegant idea that was later used to solve the approximate version \[77, 52\], and that has been recently generalized as the labeled product \[78\]. The lower bound of Equi et al. \[34\] shows that the problem remains quadratic (under SETH) even if the problem is restricted to deterministic DAGs with vertices of two kinds: indegree at most 1 and outdegree 2, and indegree 2 and outdegree at most 1 \[34\], or if restricted to undirected graphs with degree at most 2 \[34\]. Furthermore, they show how to solve the remaining cases (in/out-trees whose roots can be connected by a cycle) in linear time by an extension of KMP. Later they showed \[35, 36\] that the quadratic lower bound holds even when allowing polynomial indexing time.

An (important) special case. Gagie et al. \[42\] introduced Wheeler graphs as a generalization of prefix sortable techniques \[38, 45, 37, 22, 83\] applied to labeled graphs. On Wheeler graphs, SMLG can be solved in time \(O(m \log |E|)\) \[42\] after indexing, however, it was shown that the languages recognized by Wheeler graphs (intuitively the set of strings they encode) are very restrictive \[5, 6\]. Later, Cotumaccio and Prezza \[31\] generalized Wheeler graphs to \(p\)-sortable graphs, capturing every labeled graph by using the parameter \(p\): the minimum width of a colex relation over the vertices of the graph. On \(p\)-sortable graphs, SMLG can be solved in time \(O(mp^2 \log (pq))\) after indexing, however, the problems of deciding if a labeled graph is Wheeler or \(p\)-sortable are NP-hard \[44\]. In a recent work, Cotumaccio \[30\] defined \(q\)-sortable graphs as a relaxation of \(p\)-sortable (\(q < p\)), which can be indexed in \(O(|E|^2 + |V|^{5/2})\) time but still solve SMLG in time \(O(nq^2 \log (qq))\).

1.1 Our results

We present parameterized algorithms for SMLG on DAGs. Our parameters capture the topological structure of \(G\). These results are related to the line of research “FPT inside P” \[43, 26, 40, 61, 1, 23, 25, 24, 66, 64\] of finding parameterizations for polynomially-solvable problems.

All our results are derived from a new version of the DAG algorithm \[74\], which we present in Section 3. Our algorithm is optimized to only carry prefix-incomparable matches (Definition 3) and process them in time proportional to their size (Lemma 5 further optimized in Lemma 6 and Theorem 7). Prefix-incomparable sets suffice to capture all prefix matches

\(^2\) \(N\) is the total length of the labels in \(G\) in a more general version of the problem where vertices are labeled with strings.
of $S$ ending in a vertex $v$ (Definition 4). By noting that the size of prefix-incomparable sets is upper-bounded by the structure of $S$ (Lemma 27 in Appendix A), we obtain a parameterized algorithm (Theorem 28 in Appendix A) that beats the DAG algorithm on periodic strings.

To obtain the parameterization on the topological structure of the graph, we first study and generalize a special class of DAGs called funnels in Section 4.

**Funnels.** Funnels are DAGs whose source-to-sink paths contain a private edge that is not used by any other source-to-sink path. Although more complex that in/out-forests, their simplicity has allowed to efficiently solve problems that remain hard even when the input is restricted to DAGs, including: DAG partitioning [70], k-linkage [70], minimum flow decomposition [57, 56], a variation of network inhibition [62] and SMLG (this work). Millani et al. [70] showed that funnels can also be characterized by a partition into an in-forest plus an out-forest (the vertex partition characterization), or by the absence of certain forbidden paths (the forbidden path characterization), and propose how to find a minimal forbidden path in quadratic $O(|V|(|V|+|E|))$ time and a recognition algorithm running in $O(|V|+|E|)$ time. They used the latter to develop branching algorithms for the NP-hard problems of vertex and edge deletion distance to a funnel, obtaining a fix-parameter quadratic solution. Analogous to the minimum feedback set problem [54], the vertex (edge) deletion distance to a funnel problem asks to find the minimum number of vertices (edges) that need to be removed from a graph so that the resulting graph is a funnel.

We propose three (new) linear time recognition algorithms of funnels (Section 4.1), each based on a different characterization, improving the running time of the branching algorithm to parameterized linear time (see Appendix B). We generalize funnels to $k$-funnels by allowing private edges to be shared by at most $k$ source-to-sink paths (Definitions 14 and 15). We show how to recognize them in linear time (Lemma 16) and find the minimum $k$ for which a DAG is a $k$-funnel (Corollary 17 and Lemma 18). We then further generalize $k$-funnels to the class of DAGs $ST_k$ (Definition 20 and Lemma 21), which (unlike $k$-funnels for $k > 1$, see Figure 2) can be characterized (and efficiently recognized, see Lemma 22) by a partition into a graph of the class $S_k$ (generalization of out-forest, see Definition 19) and a graph of the class $T_k$ (generalization of in-forest, see Definition 19).

We obtain our parameterized results in Section 5 by noting that, analogous to the fact that in KMP we only need the longest prefix match, in the DAG algorithm we can bound the size of the prefix-incomparable sets by the number of paths from a source or the number of paths to a sink, $\mu_s(v)$ and $\mu_t(v)$, respectively (Lemma 23).

**Theorem 1.** Let $G = (V,E)$ be a DAG, $\Sigma$ a finite alphabet ($\sigma = |\Sigma|$), $\ell : V \rightarrow \Sigma$ a labeling function and $S \in \Sigma^m$ a string. We can decide whether $S$ has a match in $(G,\ell)$ in time $O((|V|+|E|)k + \sigma m)$, where $k = \min(\max_{v \in V} \mu_s(v), \max_{v \in V} \mu_t(v))$.

In particular, this implies linear time algorithms for out-forests and in-forests, and for every DAG in $S_k$ or $T_k$ for constant $k$. Finally, we solve the problem on DAGs in $ST_k$ (thus also in $k$-funnels), by using the vertex partition characterization of $ST_k$ (Lemma 22), solving the matches in each part separately with Theorem 1, and resolve the matches crossing from one part to the other with a precomputed data structure (Lemma 26).

**Theorem 2.** Let $G = (V,E)$ be a DAG, $\ell : V \rightarrow \Sigma$ a labeling function and $S \in \Sigma^m$ a string. We can decide whether $S$ has a match in $(G,\ell)$ in time $O((|V|+|E|)k^2 + m^2)$, where $k = \max_{v \in V}(\min(\mu_s(v), \mu_t(v)))$. 
2 Preliminaries

We work with a (directed) graph $G = (V, E)\), a function $\ell : V \to \Sigma$ labeling the vertices of $G$ with characters from a finite alphabet $\Sigma$ of size $\sigma$, and a sequence $S[1..m] \in \Sigma^m$.

Graphs. A graph $H = (V_H, E_H)$ is said to be a subgraph of $G$ if $V_H \subseteq V$ and $E_H \subseteq E$. If $V' \subseteq V$, then $G[V']$ is the subgraph induced by $V'$, defined as $G[V'] = (V', \{(u, v) \in E : u, v \in V'\})$. We denote $G^c = (V, E^c)$ to be the reverse of $G$ ($E^c = \{(u, v) \mid (u, v) \in E\}$).

For a vertex $v \in V$ we denote by $N^+_v$ ($N^-_v$) the set of in(out)-neighbors of $v$, and by $d^+_v = |N^+_v|$ ($d^-_v = |N^-_v|$) its in(out)degree. A source (sink) is a vertex with zero in(out)degree. The edge contraction of $(u, v) \in E$ is the graph operation that removes $(u, v)$ and merges $u$ and $v$. A path $P$ is a sequence $v_1, \ldots, v_{|P|}$ of different vertices of $V$ such that $(v_i, v_{i+1}) \in E$ for every $i \in \{1, \ldots, |P| - 1\}$. We say that $P$ is proper if $|P| \geq 2$, a cycle if $(v_{|P|}, v_1) \in E$, and source-to-sink if $v_1$ is a source and $v_{|P|}$ is a sink. We say that $u \in V$ reaches $v \in V$ if there is a path from $u$ to $v$. If $G$ does not have cycles it is called directed acyclic graph (DAG). A topological ordering of a DAG is a total order $v_1, \ldots, v_{|V|}$ of $V$ such that for every $(v_i, v_j) \in E$, $i < j$. It is known [53, 84] how to compute a topological ordering in $O(|V| + |E|)$ time, and we assume one ($v_1, \ldots, v_{|V|}$) is already computed if $G$ is a DAG. An out(in)-forest is a DAG such that every vertex has out(in)degree at most one, if it has a unique source (sink) it is called an out(in)-tree. The label of a path $P = v_1, \ldots, v_{|P|}$ is the sequence of the labels of its vertices, $\ell(P) = \ell(v_1) \ldots \ell(v_{|P|})$.

Strings. We say that $S$ has a match in $(G, \ell)$ if there is a path whose label is equal to $S$, every such path is an occurrence of $S$ in $(G, \ell)$. We denote $S[i..j]$ (also $S[i]$ if $i = j$, and the empty string if $j < i$) to be the substring of $S$ between position $i$ and $j$ (both inclusive), we say that it is proper if $i > 1$ or $j < m$, a prefix if $i = 1$ and a suffix if $j = m$. We denote $S^c$ to be the reverse of $S$ ($S^c[i] = S[m - i + 1]$ for $i \in \{1, \ldots, m\}$). A substring of $S$ is called a border if it is a proper prefix and a proper suffix at the same time. The failure function of $S$, $f_S : \{1, \ldots, m\} \to \{0, \ldots, m\}$ (just $f$ if $S$ is clear from the context), is such that $f_S(i)$ is the length of the longest border of $S[1..i]$. We also use $f_S$ to denote the in-tree $\{(0, \ldots, m), (i, f_S(i)) \mid i \in \{1, \ldots, m\}\}$, also known as the failure tree [46] of $S$. By definition, the lengths of all borders of $S[1..i]$ in decreasing order are $f_S(i), f_S^2(i), \ldots, 0$. The matching automaton of $S$, $A_S : \{0, \ldots, m\} \times \Sigma \to \{0, \ldots, m\}$, is such that $A_S(i, a)$ is the length of the longest border of $S[1..i] \cdot a$. It is known how to compute $f_S$ in time $O(m)$ [60] and $A_S$ in time $O(\sigma m)$ [2, 29, 81], and we assume they are already computed.

3 The DAG algorithm on prefix-incomparable matches

A key idea in our linear time parameterized algorithm is that of prefix-incomparable sets of the string $S$. We will show that one prefix-incomparable set per vertex suffices to capture all the matching information. See Figure 1 for an example of these concepts.

Definition 3 (Prefix-incomparable). Let $S \in \Sigma^m$ be a string. We say that $i, j \in \{0, \ldots, m\}, i < j$ are prefix-incomparable (for $S$) if $S[1..i]$ is not a border of $S[1..j]$. We say that $B \subseteq \{0, \ldots, m\}$ is prefix-incomparable (for $S$) if for every $i < j \in B$, $i$ and $j$ are prefix-incomparable (for $S$).

3 Our algorithms run in $O(|V| + |E|)$ time.
4 Our algorithms run in $\Omega(\sigma m)$ time.
Here we consider that the empty path always exists and its label is the empty string, thus its next in-neighbor will always be its parent. That is, we maintain a prefix-incomparable set representing the prefix matches incoming from the already processed paths.

In our algorithm we will compute for each vertex $v$ a prefix-incomparable set representing the prefix matches incoming from the already processed paths. Let $f_S$ be a labeling function, and $S \in \Sigma^m$ a string. For every $v \in V$ we define the sets:

- $B_v = \{i \in \{0, \ldots, m\} | \exists P \text{ path of } G \text{ ending in } v \text{ and } \ell(P) = S[1..i]\}$
- $PI_v \subseteq B_v$ as the unique prefix-incomparable set such that for every $i \in B_v$ there is a $j \in PI_v$ such that $i = j$ or $S[1..i]$ is a border of $S[1..j]$

**Definition 4** ($B_v, PI_v$). Let $G = (V, E)$ be a DAG, $\ell : V \rightarrow \Sigma$ a labeling function, and $S \in \Sigma^m$ a string. For every $v \in V$ we define the sets:

- $B_v = \{i \in \{0, \ldots, m\} | \exists P \text{ path of } G \text{ ending in } v \text{ and } \ell(P) = S[1..i]\}$
- $PI_v \subseteq B_v$ as the unique prefix-incomparable set such that for every $i \in B_v$ there is a $j \in PI_v$ such that $i = j$ or $S[1..i]$ is a border of $S[1..j]$

**Lemma 5.** Let $G = (V, E)$ be a DAG, $v \in V$, $S \in \Sigma^m$ a string, $f_S$ its failure tree and $A_S$ its matching automaton. We can compute $PI_v$ from $PI_u$ for every $u \in N^-_v$ in time $O(w^2 \cdot d^-_v)$ or in time $O\left(k^2 \cdot \sum_{u \in N^-_v} |PI_u|\right)$, after $O(m)$ preprocessing time, where $w = |\{i \in \{0, \ldots, m\} | \exists j, f_S(j) = i\}|$.

**Proof.** We precompute constant time lowest common ancestor (LCA)$^6$ queries [3] of $f_S$ in $O(m)$ time [41, 79, 20, 17, 18, 7, 39]. Note that with this structure we can check whether $i < j$ are prefix-incomparable in constant time (LCA$(i, j) < i$).

If $v$ is a source we have that either $B_v = PI_v = \{0\}$ if $\ell(v) \neq S[1]$ or $PI_v = \{1\}$ if $\ell(v) = S[1]$, otherwise we proceed as follows. To obtain the $O(k^2)$ time, we first append all the elements of every $PI_u$ for $u \in N^-_v$ into a list $L$ (of size $k_v$), then we replace every $i \in L$ by $A_S(i, \ell(v))$, and finally we check (at most) every pair $i < j$ of elements of $L$ and test (in constant time) if they are prefix-incomparable, if they are not we remove $i$ from the list. After these $O(|L|^2) = O(k^2)$ tests $L = PI_v$.

To obtain the $O(w^2 \cdot d^-_v)$ time, we process the in-neighbors of $v$ one by one and maintain a prefix-incomparable set representing the prefix matches incoming from the already processed in-neighbors. That is, we maintain a prefix-incomparable set $PI'$, and when we process the next in-neighbor $u \in N^-_v$ we append all elements of $PI'$ and $\{A_S(i, \ell(v)) | i \in PI_u\}$ into a

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$^5$ Here we consider that the empty path always exists and its label is the empty string, thus $0 \in B_v$.

$^6$ LCA$(i, j)$ returns the lowest node in the tree $f_S$ that is ancestor of both $i$ and $j$. 

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**Figure 1** A string $S = abaababaaba$ and its failure tree $f_S$, with $w = |\{i \in \{0, \ldots, 11\} | \exists j, f_S(j) = i\}| = 5$. On the string it is shown in segmented (blue) and solid (red) lines that prefix $S[1..3] = aba$ is a border of prefix $S[1..8] = abaababa$, which can also be seen in the tree since 3 is an ancestor (parent in this case) of 8. In the tree two sets are shown $B_1 = \{1, 8\}$ in segmented (blue) circles, which is prefix-comparable, and $B_2 = \{2, 3\}$ in solid (red) circles, which is prefix-incomparable. If $B_1 \cup B_2 \cup \{0\} = B_v$ for some $v \in V$, then $PI_v = \{2, 8\}$. 

![Diagram](https://via.placeholder.com/150)
list $\mathcal{L}'$ of size $O(w)$ (by Lemma 27), then we use the same quadratic procedure applied on $\mathcal{L}$ in time $O(w^2)$ to obtain the new $PI'$. After processing all in-neighbors in time $O(d_w^{-} \cdot w^2)$, we have $PI' = PI_v$. Next, we show the correctness of both procedures.

Let $PI'_v$ be the result after applying one of the procedures explained before (that is the final state of $PI'$ or $\mathcal{L}$), by construction $PI'_v$ is prefix-incomparable. Now, consider $i \in B_v$ and a path $P$ ending in $v$ with $\ell(P) = S[1..i]$. If $P$ is of length zero (only one vertex), then $i = 1$ and $\ell(P) = \ell(v) = S[1]$. Consider any $u \in N_w^-$, and any $j \in PI_u$ (there is at least one since $0 \in B_u$), this value is mapped to $A_S(j, \ell(v)) = A_S(j, S[1]) = j'$. By definition of the matching automaton, $j'$ is the longest border of $S[1..j] \cdot S[1]$, thus $S[1]$ is a border of $S[1..j']$ or $j' = 1$, in both procedures $j'$ can only be removed from $PI'_v$ if a longer prefix contains $S[1..j']$ as a border and also $S[1]$. If $P$ is a proper path and $i > 1$, consider the second to last vertex $u$ of $P$. Note that $i - 1 \in B_u$, and thus there is $j \in PI_u$, such that $S[1..i - 1]$ is a border of $S[1..j]$, this value is mapped to $j' = A_S(j, \ell(v)) = A_S(j, S[i])$, which is the length of the longest border of $S[1..j] \cdot S[i]$, but since $S[1..i]$ is also a border of $S[1..j] \cdot S[i]$, then $S[1..i]$ is also a border of $S[1..j']$ or $j' = i$. Again, in both procedures $j'$ can only be removed from $PI'_v$ if a longer prefix contains $S[1..j']$ as a border and also $S[1..i]$. Finally, we note that $PI'_v \subseteq B_v$ since every $i \in PI'_v$ corresponds to a match of $S[1..i]$ by construction. ▲

We improve the dependency on $w$ and $k_v$ by replacing the quadratic comparison by sorting plus a linear time algorithm on the balanced parenthesis representation [51, 71] of $f_S$.

\begin{lemma}
We can obtain Lemma 5 in time $O(sort(w, m) \cdot d_w^-)$ or in time $O(sort(k_v, m))$, where $sort(n, p)$ is the time spent by an algorithm sorting $n$ integers in the range $\{0, \ldots, p\}$.
\end{lemma}

\begin{proof}
We compute the balanced parenthesis (BP) [51, 71] representation of the topology of $f_S$, that is, we traverse $f_S$ from the root in preorder, appending an open parenthesis when we first arrive at a vertex, and a closing one when we leave its subtree. As a result we obtain a balanced parenthesis sequence of length $2(m + 1)$, where every vertex $i \in f_S$ is mapped to its open parenthesis position $open[i]$ and to its close parenthesis position $close[i]$, which can be computed and stored at preprocessing time. Note that in this representation, $i$ is ancestor of $j$ (and thus prefix-comparable) if and only if $open[i] \leq open[j] \leq close[i]$. As such, if we have a list of $O(k_v)$ (or $O(w)$) ($\mathcal{L}$ and $\mathcal{L}'$ from Lemma 5) values, we can compute the corresponding prefix-incomparable sets as follows.

First, we sort the list by increasing open value, this can be done in $O(sort(k_v, m))$ (or $O(sort(w, m))$, since this sorting is equivalent to sort by increasing $open/2 \in \{0, \ldots, m\}$ value. Then, we process the list in the sorted order, if two consecutive values $i$ and $j$ in the order are prefix-comparable (that is, if $open[j] \leq close[i]$) then we remove $i$ and continue to the next value $j$. At the end of this $O(k_v)$ (or $O(w)$) time processing we obtain the desired prefix-incomparable set. ▲

If we use techniques for integer sorting [87, 90, 59] we can get $O(k_v \log \log m)$ (or $O(w \log \log m)$ time for sorting, however introducing $m$ into the running time. We can solve this issue by using more advanced techniques [48, 9, 47] obtaining an $O(k_v \log \log k_v)$ (or $O(w \log \log w)$) time for sorting. However, we show that by using the suffix-tree [89, 68, 85] of $S'$ we can obtain a linear dependency on $w$ and $k_v$.

\begin{theorem}
We can obtain the result of Lemma 5 in time $O(w \cdot d_w^-)$ or in time $O(k_v)$.
\end{theorem}

\begin{proof}
We reuse the procedure of Lemma 6 but this time on the BP representation of the topology of the suffix-tree $T_r$ of $S'$, which has $O(m)$ vertices and can be built in $O(m)$ time [89, 68, 85]. Note that every suffix represented in $T_r$ corresponds to a prefix of $S$ (spelled
in the reverse direction). Moreover, \( i \leq j \) are prefix-comparable if and only if the vertex representing \( i \) in \( T_r \) is an ancestor of \( T_r[j] \), the same property as in \( f_S \). Furthermore, if \( B \) is prefix-incomparable and \( A(j, a) = j + 1 \), then the positions of the vertices in \( A(B, a) \) in \( T_r \) follow the same order as the ones in \( B \), since the suffix-tree is lexicographically sorted.

Now, we show how to obtain the prefix-incomparable set representing \( A(PI_u, \ell(v)) \) in \(|PI_u|\) time assuming that \( PI_u \) is sorted by increasing (open) position in \( T_r \).

We first separate \( PI_u \) into the elements \( i \in M \) with \( S[i + 1] = \ell(v) \) and \( i \in E \) with \( S[i + 1] \neq \ell(v) \) (in the same relative order as in \( PI_u \), which is supposed to be in increasing order). Since \( M \) is prefix-incomparable the positions of the vertices in \( A(M, \ell(v)) \) in \( T_r \) follow the same order as the ones in \( M \). We then obtain the list \( E_u \) by applying \( T_r[A(i, \ell(v)) - 1] \) for every \( i \in E \) (if \( A(i, \ell(v)) = 0 \) we do not process \( i \)), and then for any pair of consecutive elements \( x \) before \( y \) in \( E_u \) such that \( y \leq x \) we remove \( y \) from \( E_u \), and repeat this until no further such inversion remains, thus obtaining an increasing list in \( E_u \) representing vertices in \( T_r \). Next, since \( E_u \) is sorted we can obtain the list \( PI_E \) of prefix-incomparable elements representing \( E_u \), and finally apply \( A(PI_E, \ell(v)) \) (which also follows an increasing order in \( T_r \), merge it with \( A(M, \ell(v)) \), and compute the prefix-comparable elements of this merge.

The correctness of the previous procedure follows by the fact that if there is an inversion \( y < x \) in \( E_u \), then the prefix \( A(j, \ell(v)) - 1 \) represented by \( y \) in \( T_r \) is a border of the prefix \( A(i, \ell(v)) - 1 \) represented by \( x \) (and thus is safe to remove \( y \)). For this, first note that \( i \) appears before \( j \) in \( E \), then \( S[i..1] <_{tex} S[j..1] \), and since \( i \) is prefix-incomparable with \( j \) there is a \( k \geq 1 \) such that \( S[i..i + k - 1] = S[j..j + k - 1] \) and \( S[i + k] <_{tex} S[j + k] \). Then, since \( y \) appears before \( x \) in \( E_u \), then \( S[A(j, \ell(v)) - 1..1] <_{tex} S[A(i, \ell(v)) - 1..1] \), but since \( A(i, \ell(v)) - 1 \) is a border of \( i \) and \( A(j, \ell(v)) - 1 \) is a border of \( j \), \( S[A(j, \ell(v)) - 1..1] \) must be a prefix of \( S[A(i, \ell(v)) - 1..1] \), and thus \( S[1..A(j, \ell(v)) - 1] \) is a border of \( S[1..A(i, \ell(v)) - 1] \).

The corollary is obtained by maintaining the \( PI_u \) sets sorted by position in \( T_r \), and noting that the previous procedure runs in linear \( O(|PI_u|) \) time.

In Appendix A we show how to use Theorem 7 to derive a parameterized algorithm using parameter \( w = \{|i \in [0, \ldots, m] | \exists j, f_S(j) = i\} \), improving on the classical DAG algorithm when \( S \) is a periodic string. Next, we will present our results on recognizing funnels and their generalization (Section 4), and how to use these classes of graphs and Theorem 7 to obtain parameterized algorithms using parameters related to the topology of the DAG (Section 5).

## 4 Funnel and beyond

Recall that funnels are DAGs whose source-to-sink paths have at least one private edge\(^7\), that is, an edge used by only one source-to-sink path. More formally,

\begin{definition}[Private edge] Let \( G = (V, E) \) be a DAG and \( \mathcal{P} \) the set of source-to-sink paths of \( G \). We say that \( e \in E \) is private if \( \mu(e) := \{|P \in \mathcal{P} | e \in P\} = 1 \). If \( \mu(e) > 1 \), we say that \( e \) is shared. \end{definition}

\begin{definition}[Funnel] Let \( G = (V, E) \) be a DAG and \( \mathcal{P} \) the set of source-to-sink paths of \( G \). We say that \( G \) is a funnel if for every \( P \in \mathcal{P} \) there exists \( e \in P \) such that \( e \) is private. \end{definition}

Millani et al. [70] showed two other characterizations of funnels.

\(^7\) For the sake of simplicity, we assume that there are no isolated vertices, thus any source-to-sink path has at least one edge.
**Theorem 10** ([70]). Let $G = (V, E)$ be a DAG. The following are equivalent:
1. $G$ is a funnel
2. There exists a partition $V = V_1 \cup V_2$ such that $G[V_1]$ is an out-forest, $G[V_2]$ is an in-forest and there are no edges from $V_2$ to $V_1$
3. There is no path $P$ such that its first vertex has more than one in-neighbor (a merging vertex) and its last vertex more than one out-neighbor (a forking vertex). Such a path is called forbidden.

They also gave an $O(|V| + |E|)$ time algorithm to recognize whether a DAG $G$ is a funnel, and an $O(|V|(|V| + |E|))$ time algorithm to find a minimal forbidden path in a general graph, that is, a forbidden path that is not contained in another forbidden path.

### 4.1 Three (new) linear time recognition algorithms

We first show how to find a minimal forbidden path in time $O(|V| + |E|)$ in general graphs, improving on the quadratic algorithm of Millani et al. [70].

**Lemma 11.** Let $G = (V, E)$ be a graph. In $O(|V| + |E|)$ time, we can decide if $G$ contains a forbidden path, and if one exists we report a minimal forbidden path.

**Proof.** In the bioinformatics community minimal forbidden paths are a subset of unitigs and it is well known how to compute them in $O(|V| + |E|)$ time (see e.g. [55, 50, 58, 69]), here we include a simple algorithm for completeness. We first compute the indegree and outdegree of each vertex and check whether there exists a forbidden path of length zero or one, all in $O(|V| + |E|)$ time, in the process we also mark all vertices except the ones with unit indegree and outdegree. If no path is found we iterate over the vertices one last time. If the current vertex is not marked we extend it back and forth to the closest marked vertices and mark the vertices in these extensions, finally we check whether the first vertex is merging and the last forking. This last step takes $O(|V|)$ time in total.

Lemma 11 provides our first linear time recognition algorithm and, as opposed to the algorithm of Millani et al. [70], it also reports a minimal forbidden path given a general graph. Moreover, in Appendix B, we show that Lemma 11 provides a linear time parameterized algorithm for the NP-hard (and inapproximable) problem of deletion distance of a general graph to a funnel [63, 70]. Millani et al. [70] solved this problem in (parameterized) quadratic time and in (parameterized) linear time only if the input graph is a DAG.

Next, we show another linear time recognition algorithm, which additionally finds the partition $V = V_1 \cup V_2$ from Theorem 10. Finding such a partition will be essential for our solution to SMLG. From now we will assume that the input graph is a DAG since this condition can be checked in linear time [53, 84].

**Lemma 12.** Let $G = (V, E)$ be a DAG. We can decide in $O(|V| + |E|)$ time whether $G$ is a funnel. Additionally, if $G$ is a funnel, the algorithm reports a partition $V = V_1 \cup V_2$ such that $G[V_1]$ is an out-forest, $G[V_2]$ is an in-forest and there are no edges from $V_2$ to $V_1$.

**Proof.** We start a special BFS traversal from all the source vertices of $G$. The traversal only adds vertices to the BFS queue if they have not been previously visited (as a typical BFS traversal) and if its indegree is at most one. After the search we define the partition $V_1$ as the set of vertices visited during the traversal and $V_2 = V \setminus V_1$. Finally, we report the previous partition if there are no edges from $V_2$ to $V_1$, and if every vertex of $V_2$ has outdegree at most one. All these steps run in time $O(|V| + |E|)$. 

Note that if the algorithm reports a partition, then this satisfies the required conditions to be a funnel \((G[V_1]\) is an out-forest since every vertex visited in the traversal has indegree at most one). Moreover, if \(G\) is a funnel, we prove that \(V_2\) is an in-forest and that there are no edges from \(V_2\) to \(V_1\). For the first, suppose by contradiction that there is a vertex \(v \in V_2\) with \(d^+_v > 1\), since every vertex is reached by some source in a DAG then there is a \(u \in V_2\) with \(d^-_u > 1\) (a vertex that was not added to the BFS queue) that reaches \(v\), implying the existence of a forbidden path in \(G\), a contradiction. Finally, there cannot be edges from \(V_2\) to \(V_1\) since the indegree (in \(G\)) of vertices of \(V_1\) is at most one and its unique (if any) in-neighbor is also in \(V_1\) by construction.

Next, we present another characterization of funnels based on the structure of private/shared edges of the graph, which can be easily obtained by manipulating the original Definition 9.

\textbf{Definition 13 (Funnel).} Let \(G = (V, E)\) be a DAG. We say that \(G\) is a funnel if there is no source-to-sink path using only shared edges.

As such, another approach to decide whether a DAG \(G\) is a funnel is to compute \(\mu(e)\) for every \(e \in E\) and then perform a traversal that only uses shared edges. Computing the number of source-to-sink paths containing \(e\), that is \(\mu(e)\), can be done by multiplying the number of source-to-\(e\) paths, \(\mu_s(e)\), by the number of \(e\)-to-sink paths, \(\mu_t(e)\), each of which can be computed in \(O(|V| + |E|)\) time for all edges. The solution consists of a dynamic program on a topological order (and reverse topological order) of \(G\) with the following recurrences.

\[
\begin{align*}
\mu_s(u) &= \mathbb{1}_{d^-_u = 0} + \sum_{u' \in N^-_u} \mu_s(u') \\
\mu_t(v) &= \mathbb{1}_{d^+_v = 0} + \sum_{v' \in N^+_v} \mu_t(v') \\
\mu((u, v)) &= \mu_s(u) \cdot \mu_t(v)
\end{align*}
\]

Where \(\mathbb{1}_A\) is the characteristic function evaluating to 1 if \(A\) is true and to 0 otherwise. It is simple to observe that the previous dynamic programs can be computed in time \(O(|V| + |E|)\) each and that for every \(e = (u, v) \in E\), \(\mu_s(e) = \mu_s(u)\) and \(\mu_t(e) = \mu_t(v)\). By simplicity, in the following we will use \(\mu_s(e)\) and \(\mu_s(u)\) (also \(\mu_t(e)\) and \(\mu_t(v)\)) interchangeably. The previous algorithm assumes constant time arithmetic operations on numbers up to \(\max_{e \in E} \mu(e)\), which can be \(O(2^{|V|})\). To avoid this issue, we note that it is not necessary to compute \(\mu(e)\), but only to verify that \(\mu(e) > 1\). As such, we can recognize shared edges as soon as we identify that \(\mu(e) > 1\), that is whenever \(\mu_s(e)\) or \(\mu_t(e)\) is greater than one in their respective computation. A formal description of this algorithm can be found in Lemma 16.

\subsection{4.2 Generalizations of funnels}

To generalize funnels we will allow source-to-sink paths to use only shared edges, but require to have at least one edge shared by at most \(k\) different source-to-sink paths.

\textbf{Definition 14 \((k\text{-private edge})\).} Let \(G = (V, E)\) be a DAG. We say that \(e \in E\) is \(k\)-private if \(\mu(e) \leq k\). If \(\mu(e) > k\) we say that \(e\) is \(k\)-shared.

\textbf{Definition 15 \((k\text{-funnel})\).} Let \(G = (V, E)\) be a DAG. We say that \(G\) is a \(k\)-funnel if there is no source-to-sink path using only \(k\)-shared edges.
The next algorithm is a generalization of the last algorithm in Section 4.1 to decide if a DAG is a $k$-funnel. It assumes constant time arithmetic operations on numbers up to $k$.

**Lemma 16.** We can decide if a DAG $G = (V, E)$ is a $k$-funnel in $O(|V| + |E|)$ time, assuming constant time arithmetic operations on numbers up to $\Theta(k)$.

**Proof.** We process the vertices in a topological ordering and use Equation (1) to compute $\mu_s(e)$ in one pass, $\mu_t(e)$ in another pass and $\mu(e)$ in a final pass. To avoid arithmetic operations with numbers greater than $k$, we mark the edges having $\mu_s$ and $\mu_t$ greater than $k$ as $k$-shared during the computations of $\mu_s, \mu_t$. Note that if $\mu_s(e) > k$ or $\mu_t(e) > k$ then $\mu(e) > k$. As such, before computing $\mu_s(e)$ ($\mu_t(e)$) we check if some of the edges from (to) the in(out)-neighbors is marked as $k$-shared. If that is the case we do not compute $\mu_s(e)$ ($\mu_t(e)$) and instead mark $e$ as $k$-shared, otherwise we compute the respective sum of Equation (1), and if at some point the cumulative sum exceeds $k$ we stop the computation and mark $e$ as $k$-shared. Finally, we find all $k$-shared edges as the marked plus the unmarked with $\mu(e) = \mu_s(e) \cdot \mu_t(e) > k$, perform a traversal only using $k$-shared edges, and report that $G$ is not a $k$-funnel if there is a source-to-sink path using only $k$-shared edges in time $O(|V| + |E|)$.

We can use the previous result and exponential search [19, 15] to find the minimum $k$ such that a DAG is a $k$-funnel.

**Corollary 17.** Let $G = (V, E)$ be a DAG. We can find the minimum $k$ such that $G$ is a $k$-funnel in $O((|V| + |E|) \log k)$ time, assuming constant time arithmetic operations on numbers up to $\Theta(k)$.

Assuming constant time arithmetic operations on numbers up to $\max_{e \in E} \mu(e)$ the problem is solvable in linear time by noting that the answer is equal to the weight of a widest path.

**Lemma 18.** Let $G = (V, E)$ be a DAG. We can find the minimum $k$ such that $G$ is a $k$-funnel in $O(|V| + |E|) \log k)$ time, assuming constant time arithmetic operations on numbers up to $\Theta(\max_{e \in E} \mu(e))$.

**Proof.** We compute $\mu(e)$ for every $e \in E$ by using the dynamic programming algorithm specified by Equation (1) on a topological ordering of $G$. Since constant time arithmetic operations are assumed for numbers up to $\max_{e \in E} \mu(e)$, the previous computation takes linear time. Then, we compute the weight of a source-to-sink path $P$ maximizing $\min_{e \in P} \mu(e)$, and report this value. This problem is known as the widest path problem [76, 82, 65, 86, 80] and it can be solved in linear time on DAGs [88, 49] by a dynamic program on a topological order of the graph. By completeness, we show a dynamic programming recurrence to compute $W[e]$, the weight of a source-to-e path $P$ maximizing $\min_{e' \in P} \mu(e')$.

$$W[e = (u, v)] = \mu(e) \cdot \mathbb{1}_{d^-_e = 0} + \sum_{w' \in N^-_e} \min(W[(u', u), \mu((u', u))])$$

Finally, note that if we denote $w$ to the weight of a widest path, then there is a source-to-sink path using only $w - 1$-shared edges, $G$ is not $w - 1$-funnel. Moreover, there cannot be a source-to-sink path using only $w$-shared edges, since such a path would contradict $w$ being the weight of a widest path. As such, $w$ is the minimum $k$ such that $G$ is $k$-funnel.

We now define three classes of DAGs closely related to $k$-funnels.

**Definition 19.** We say that a DAG $G = (V, E)$ belongs to the class $S_k$ ($T_k$) if for every $v \in V$, $\mu_s(v)$ ($\mu_t(v)$) $\leq k$. 
The red label next to each edge corresponds to \( \min(\mu_s(v), \mu_t(v)) \), since the maximum of these labels is \( k \) the graph belongs to \( ST_k \). The red label next to each edge \( e \) corresponds to \( \mu(e) \), since there is a source-to-sink path with no \( k \)-private edge the graph is not a \( k \)-funnel.

\[ \text{Definition 20.} \text{ We say that a DAG } G = (V, E) \text{ belongs to the class } ST_k \text{ if for every } v \in V, \mu_s(v) \leq k \text{ or } \mu_t(v) \leq k. \]

\[ \text{Lemma 21.} \text{ } S_k, T_k \subseteq k\text{-funnels } \subseteq ST_k. \]

\textbf{Proof.} We first prove that \( S_k, T_k \subseteq k\text{-funnels} \). Consider \( G \in S_k \) \( (G \in T_k) \), and take any source-to-sink path \( P \) of \( G \). Let \( (u, v) \) be the last (first) edge of \( P \), then by Equation (1) \( \mu((u, v)) = \mu_s(u) \cdot \mu_t(v) \), but since \( \mu_t(v) = 1 \) \((v) \text{ is a sink) and } \mu_s(u) \leq k \) \((G \in S_k) \) analogously, \( \mu_s(u) = 1 \text{ and } \mu_t(v) \leq k \), then \( \mu((u, v)) \leq k \), and thus \( (u, v) \) is a \( k \)-private edge. To prove that \( k\)-funnels \( \subseteq ST_k \), suppose that \( G \) is a \( k \)-funnel, and by contradiction that there exists \( v \in V \) with \( \mu_s(v), \mu_t(v) > k \). Consider any source-to-sink path \( P \) using \( v \). Now, let \( (u, w) \) be any edge in \( P \) before (after) \( v \), then \( \mu_t(w) \geq \mu_t(v) > k \) \((\mu_s(u) \geq \mu_t(v) > k) \), and thus \( \mu((u, w)) = \mu_s(u) \cdot \mu_t(w) > k \). As such, \( P \) does not have a \( k \)-private edge, a contradiction. \( \square \)

For \( k = 1 \), \( S_1 \) describes out Forests and \( T_1 \) in-forests, thus being more restrictive than funnels. Moreover, we note that the in(out)-star of \( k + 2 \) vertices, that is \( k + 1 \) vertices pointing to a sink (pointed from a source), \( \not\in \mathcal{S}_k \) \((\mathcal{T}_k) \), but this graph is a funnel. On the other hand, from the vertex partition characterization of funnels (Theorem 10 \([70]\)) we have that \( ST_1 = \{1\}\text{-funnels} \). However, for \( k \), the containment \( k\text{-funnels } \subseteq ST_k \) is strict (Figure 2).

By noting that the minimum \( k \) such that a DAG is in \( S_k \), \( T_k \) and \( ST_k \) is \( \max_{v \in V} \mu_s(v) \), \( \max_{v \in V} \mu_t(v) \) and \( \max_{v \in V} \min(\mu_s(v), \mu_t(v)) \), respectively, we obtain the same results as in Lemmas 16 and 18 and Corollary 17 (with analogous assumptions on the cost of arithmetic operations) for recognition of \( S_k \), \( T_k \) and \( ST_k \).

Next, we prove that although the vertex partition characterization of funnels does not generalize to \( k \)-funnels, it does for the class \( ST_k \) and it can be found efficiently.

\[ \text{Lemma 22.} \text{ Let } G = (V, E) \in ST_k \text{ and } k \text{ given as inputs. We can find, in } O(|V| + |E|) \text{ time, a partition } V = V_1 \cup V_2 \text{ such that } G[V_1] \in S_k, G[V_2] \in T_k \text{ and there are no edges from } V_2 \text{ to } V_1 \text{. Moreover, if such a partition of a DAG } G \text{ exists, then } G \in ST_k. \]

\textbf{Proof.} We set \( V_1 = \{ v \in V \mid \mu_s(v) \leq k \} \) and \( V_2 = V \setminus V_1 \). Note that finding \( V_1 \) takes linear time, since we can apply the algorithm described in Lemma 16 to compute the \( \mu_s \) values (or decide that they are more than \( k \))\(^8\). By construction we know that every \( v \in V_1 \) has

\(^8\) Recall that this procedure assumes constant time arithmetic operations of numbers up to \( \Theta(k) \).
\( \mu_s(v) \leq k \), and since \( G \in ST_k \) also every \( v \in V_2 \) has \( \mu_t(v) \leq k \), thus \( G[V_1] \in S_k \), \( G[V_2] \in T_k \).

Suppose by contradiction that there exists \( e = (u, v) \in E \cap (V_2 \times V_1) \). As such, \( \mu_s(u) > k \), but since \( \mu_s(u) \leq \mu_s(v) \), then \( \mu_s(v) > k \), a contradiction. Finally, if such a partition exists then \( \mu_s(v) \leq k \) for every \( v \in V_1 \) and \( \mu_t(v) \leq k \) for every \( v \in V_2 \), and thus \( G \in ST_k \).

### 5. Parameterized algorithms: The DAG

The main idea to get the parameterized algorithms in this section is to bound the size of the \( PI_v \) sets by a topological graph parameter and use Lemma 5 and Theorem 7 to obtain a parameterized solution. As in the KMP algorithm [60] only one prefix-incomparable value suffices (the longest prefix match until that point), we show that \( \mu_s(v) \) prefix-incomparable values suffice to capture the prefix matches up to \( v \).

**Lemma 23.** Let \( G = (V, E) \) be a DAG, \( v \in V \), \( \mathcal{P}_v \) the set of source-to-\( v \) paths, \( \ell : V \to \Sigma \) a labeling function, \( S \in \Sigma^m \) a string, and \( PI_v \) as in Definition 4. Then, \( |PI_v| \leq \mu_s(v) \).

**Proof.** Since any path ending in \( v \) is the suffix of a source-to-\( v \) path we can write \( B_v \) as:

\[
B_v = \bigcup_{P_{sv} \in \mathcal{P}_v} B_{P_{sv}} : = \{ i \in \{0, \ldots, m \} \mid \exists \text{ suffix of } P_{sv}, \ell(P) = S[1..i] \}
\]

However, for every pair of values \( i < j \in B_{P_{sv}} \), \( S[1..i] \) is a border of \( S[1..j] \) (it is a suffix since they are both suffixes of \( \ell(P_{sv}) \)). As such, at most one value of \( B_{P_{sv}} \) appears in \( PI_v \), and then \( |PI_v| \leq |\mathcal{P}_sv| = \mu_s(v) \).

This result directly implies a parameterized string matching algorithm to DAGs in \( S_k \).

**Lemma 24.** Let \( G = (V, E) \in S_k \), \( \ell : V \to \Sigma \) a labeling function and \( S \in \Sigma^m \) a string. We can decide whether \( S \) has a match in \( (G, \ell) \) in time \( O(|V|k + |E| + \sigma m) \).

**Proof.** We compute the matching automaton \( A_S \) in \( O(\sigma m) \) time. Then, we process the vertices in topological order, and for each vertex \( v \) we compute \( PI_v \), the unique prefix-incomparable set representing \( B_v \) (all prefix matches of \( S \) with paths ending in \( v \)). We proceed according to Lemma 5 and Theorem 7 in \( O(m) \) preprocessing time plus \( O(k_v) \) time per vertex. There is a match of \( S \) in \( (G, \ell) \) if and only if any \( PI_v \) contains \( m \). The claimed running time follows since \( k_v = \sum_{u \in N_v} |PI_u| \leq \sum_{u \in N_v} \mu_s(u) \leq \mu_s(v) \leq k \), by Lemma 23, Equation (1) and since \( G \in S_k \).

A simple but interesting property about string matching to graphs is that we obtain the same problem by reversing the input (both the graph and the string), that is, \( S \) has a match in \( (G, \ell) \) if and only if \( S^r \) has a match in \( G^r, \ell \). This fact, plus noting that \( G \in S_k \) if and only if \( G^r \in T_k \) gives the following corollary of Lemma 24.

**Corollary 25.** Let \( G = (V, E) \in T_k \), \( \ell : V \to \Sigma \) a labeling function and \( S \in \Sigma^m \) a string. We can decide whether \( S \) has a match in \( (G, \ell) \) in time \( O(|V|k + |E| + \sigma m) \).

With these two results and the fact that we can compute the minimum \( k \) such that a DAG is in \( S_k \), \( T_k \) in time \( O(|V| + |E|) \log k \) (see Corollary 17) we obtain our first algorithm parameterized by the topology of the DAG.

**Theorem 1.** Let \( G = (V, E) \) be a DAG, \( \Sigma \) a finite alphabet \( (\sigma = |\Sigma|) \), \( \ell : V \to \Sigma \) a labeling function and \( S \in \Sigma^m \) a string. We can decide whether \( S \) has a match in \( (G, \ell) \) in time \( O(|V| + |E|)k + \sigma m) \), where \( k = \min \{ \max_{v \in V} \mu_s(v), \max_{v \in V} \mu_t(v) \} \).
Our final result is a parameterized algorithm for DAGs in \( ST_k \) (in particular for \( k \)-funnels). We note that the algorithm of Corollary 25 computes \( PI_v \) for \( S^v \) for every vertex in \( G^* \). Recall that \( PI_v \) represents all the prefix matches of \( S^v \) with paths ending in \( v \) in \( G^* \). In other words, it represents all suffix matches of \( S \) with paths starting in \( v \) in \( G \). For clarity, let us call this set \( SI_v \). The main idea of the algorithm for \( ST_k \) is to use Lemma 22 to find a partitioning \( V = V_1 \cup V_2 \) into \( S_k \) and \( T_k \), use Lemma 24 and Corollary 25 to search for matches within each part and also to compute \( PI_v \) for every \( v \in V_1 \) and \( SI_v \) for every \( v \in V_2 \), and finally, to find matches using the edges from \( V_1 \) to \( V_2 \). The last ingredient of our algorithm consists of preprocessing the answers to the last type of matches.

▶ Lemma 26. Let \( G = (V,E) \) a DAG, \( (u,v) \in E \), \( \ell : V \to \Sigma \) a labeling function, \( S \in \Sigma^m \) a string and \( PI_u \) and \( SI_v \), as in Definition 4. We can decide if there is a match of \( S \) in \( (G,\ell) \) using \((u,v)\) in \( O(|PI_u| \cdot |SI_v|) \) time, after \( O(m^2) \) preprocessing time.

Proof. We precompute a boolean table \( PS \) of \( m \times m \) entries, such that \( PS[i,j] \) is true if there is a length \( i' \) of a (non-empty) border of \( S[1..i] \) (or \( i' = i \)) and a length \( j' \) of a (non-empty) border of \( S[m-j+1..m] \) (or \( j' = j \)) such that \( i' + j' = m \), and false otherwise. This table can be computed by dynamic programming in \( O(m^2) \) time as follows.

\[
PS[i,j] = \begin{cases} 
\text{false} & \text{if } i + j < m \lor i = 0 \lor j = 0 \\
i + j = m \lor PS[i,f_S(j)] \lor PS[f_S(i),j] & \text{otherwise}
\end{cases}
\]

We then use this table to test every \( PS[i,j] \) with \( i \in PI_u \) and \( j \in SI_v \) and report a match if any of these table entries is true, in total \( O(|PI_u| \cdot |SI_v|) \) time.

Since every match of \( S \) in \( (u,v) \) must match a prefix \( S[1..i] \) with a path ending in \( u \) and a suffix \( S[i+1..m] \) with a path starting in \( v \), the previous procedure finds it (if any).

▶ Theorem 2. Let \( G = (V,E) \) be a DAG, \( \ell : V \to \Sigma \) a labeling function and \( S \in \Sigma^m \) a string. We can decide whether \( S \) has a match in \( (G,\ell) \) in time \( O((|V| + |E|)k^2 + m^2) \), where \( k = \max_{v \in V}(\min(\mu_s(v), \mu_t(v))) \).

Proof. We first compute the minimum \( k \) such that the input DAG is in \( ST_k \) in time \( O((|V| + |E|) \log k) \) (see Corollary 17). Then, we obtain the partition of \( G \) into \( G[|V_1| \in S_k \text{ or } G[|V_2| \in T_k \text{ and no edges from } V_2 \text{ to } V_1 \). We then search matches within \( G[|V_1| \text{ and } G[|V_2| \text{ in time } O(|V|k + |E| + \sigma m) \) (Lemma 24 and Corollary 25) and we also keep \( PI_u \) for every \( u \in V_1 \) and \( SI_v \) for every \( v \in V_2 \). Finally, we process the matches using the edges \( (u,v) \) with \( u \in V_1, v \in V_2 \) in total \( O(|E|k^2 + m^2) \) time (Lemma 26) since \( O(|PI_u| \cdot |SI_v|) = O(k^2) \).

6 Conclusions

In this paper we introduced the first parameterized algorithms for matching a string to a labeled DAG, a problem known (under SETH) to be quadratic even for a very special type of DAGs. Our parameters depend on the structure of the input DAG.

We derived our results from a generalization of KMP to DAGs using prefix-incomparable matches, which allowed us to bound the running time to parameterized linear. Further improvements on the running time of our algorithms remain open: is it possible to get rid of the automaton? or to combine prefix-incomparable and suffix-incomparable matches in better than quadratic (either in the size of the sets or the string)? (e.g. with a different tradeoff between query and construction time of the data structure answering these queries) and is there a (conditional) lower bound to combine these incomparable sets? (see e.g. [21]). Another
interesting question with practical importance is whether our parameterized approach can be extended to string labeled graphs with (unparameterized) linear time in the total length of the strings or extended to counting and reporting algorithms in linear time in the number of occurrences.

We also presented novel algorithmic results on funnels as well as generalizations of them. These include linear time recognition algorithms for their different characterizations, which we showed useful for the string matching problem but hope that can also help in other graph problems. We also showed how to find the minimum $k$ for which a DAG is a $k$-funnel or $\in ST_k$ (assuming constant time arithmetic operations on numbers up to $O(k)$) using an exponential search, but it remains open whether there exists a linear time solution.

References


A simple property about prefix-incomparable sets is that their sizes are bounded by the number of prefixes that are not a border of other prefixes of the string, equivalently, the number of leaves in the failure function of the string.

Lemma 27. Let $S \in \Sigma^m$ be a string, $f_S$ its failure function/tree, and $B \subseteq \{0, \ldots, m\}$ prefix-incomparable for $S$. Then, $|B| \leq w$ such that $w$ is the number of leaves of $f_S$, equivalently $w := \{|i \in \{0, \ldots, m\} | \exists j, f_S(j) = i\}$.

Proof. First note that $i < j$ are prefix-incomparable if and only if $i$ is not ancestor of $j$ in $f_S$. Suppose by contradiction that $|B| > w$, and consider the $w$ leaf-to-root paths of $f_S$. Note that these $w$ leaf-to-root paths cover all the vertices of $f_S$. By pigeonhole principle, there must be $i < j \in B$ in the same leaf-to-root path, that is $i$ is ancestor of $j$, a contradiction.
Theorem 28. Let $G = (V, E)$ be a DAG, $\Sigma$ a finite alphabet ($\sigma = |\Sigma|$), $\ell : V \rightarrow \Sigma$ a labeling function, $S \in \Sigma^*$ a string and $f_S$ its failure function. We can decide whether $S$ has a match in $(G, \ell)$ in time $O((|V| + |E|)w + \sigma m)$, where $w = \{|i \in \{0, \ldots, m\} | \exists j, f_S(j) = i\}$.

Proof. We compute the matching automaton $A_S$ in $O(\sigma m)$ time. Then, we process the vertices in topological order, and for each vertex $v$ we compute $PI_v$, the unique prefix-incomparable set representing $B_v$ (all prefix matches of $S$ with paths ending in $v$). We proceed according to Lemma 5 and Theorem 7 in $O(m)$ preprocessing time plus $O(w \cdot d_v^2)$ time per vertex, adding up to $O(w(|V| + |E|))$ time in total. There is a match of $S$ in $(G, \ell)$ if and only if any $PI_v$ contains $m$.

We note that $w \leq m$, thus our algorithm is asymptotically as fast as the DAG algorithm, which runs in time $\Omega((|V| + |E|)m)$. However, we note that for $w = o(m)$, a (very) long prefix of $S$ must be a (highly) periodic string. To see this, consider the longest prefix $S[1..i]$ of $S$, such that there exists $j > i$ with $i = f_S(j)$. By definition, $S[1..i]$ is a border of $S[1..j]$, thus $S[k] = S[k + j - i]$ for $k \in \{1, \ldots, i\}$, that is $S[1..j]$ is a periodic string with period $j - i$. Finally, note that if $w = o(m)$, then $m - i \in o(m)$, and thus the period $j - i \in o(m)$.

B A linear time parameterized algorithm for the distance problems

Millani et al. [70] gave an $O(|V|(|V| + |E|))$ time algorithm to find a minimal forbidden path in a general graph. They used this algorithm to design branching algorithms (see e.g. [32]) for the problems of finding maximum sized sets $V' \subseteq V, d_v := |V'|$ and $E' \subseteq E, d_e := |E'|$, such that $G[V']$ and $(V, E')$ are funnels, known as vertex and edge distance to a funnel. It is known that (unless $P = NP$) there is an $\epsilon > 0$ such that there is no polynomial time $|V|^{1-\epsilon}$ approximation [63] for the vertex version nor $(1 + \epsilon)$ approximation [70] for the edge version. The authors [70] noted that if we consider a minimal forbidden path $P$ of $G$ of length $|P| > 1$, then the edges of $P$ can be contracted until $|P| = 1$ without affecting the size of the solution. Moreover, they noted that if we consider such a $P$, two in-neighbors of the first vertex and two out-neighbors of the last, then $V'$ must contain at least one of those 6 vertices and $E'$ one of those 5 edges, deriving $O(6^d |V|(|V| + |E|))$ and $O(5^d |V|(|V| + |E|))$ time branching algorithms for each problem [10] [70, Corollary 1]. The authors also developed a more involved branching algorithm, only for the edge distance problem on DAG inputs, running in time $O(3^d(|V| + |E|))$ [70, Theorem 4].

By noting that minimal forbidden paths can be further contracted to length zero (one vertex) in the vertex distance problem, and that a minimal forbidden path can be found in time $O(|V| + |E|)$ (Lemma 11) we obtain the following result.

Theorem 29. Let $G = (V, E)$ be a graph. We can compute the vertex (edge) deletion distance to a funnel in time $O(5^d(|V| + |E|))$, where $d$ is the deletion distance.

Proof. We follow the branching approach as in [70, Corollary 1], but in the case of vertex distance we further contract the forbidden paths to length 0, the correctness of this step follows by noting that any solution containing two different vertices in a forbidden path is not minimum, since we still get a funnel by removing one of them (from the solution). As such, the number of recursive calls is $\leq 5$ for both problems. Moreover, by Lemma 11, we can find a minimal forbidden path in time $O(|V| + |E|)$.
Optimal Near-Linear Space Heaviest Induced Ancestors

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Abstract

We revisit the Heaviest Induced Ancestors (HIA) problem that was introduced by Gagie, Gawrychowski, and Nekrich [CCCG 2013] and has a number of applications in string algorithms. Let $T_1$ and $T_2$ be two rooted trees whose nodes have weights that are increasing in all root-to-leaf paths, and labels on the leaves, such that no two leaves of a tree have the same label. A pair of nodes $(u, v) \in T_1 \times T_2$ is induced if and only if there is a label shared by leaf-descendants of $u$ and $v$. In an HIA query, given nodes $x \in T_1$ and $y \in T_2$, the goal is to find an induced pair of nodes $(u, v)$ of the maximum total weight such that $u$ is an ancestor of $x$ and $v$ is an ancestor of $y$.

Let $n$ be the upper bound on the sizes of the two trees. It is known that no data structure of $\tilde{O}(n)$ can answer HIA queries in $\omega(\log n / \log \log n)$ time [Charalampopoulos, Gawrychowski, Pokorski; ICALP 2020]. This (unconditional) lower bound is a polyloglog $n$ factor away from the query time of the fastest $\tilde{O}(n)$-size data structure known to date for the HIA problem [Abedin, Hooshmand, Ganguly, Thankachan; Algorithmica 2022]. In this work, we resolve the query-time complexity of the HIA problem for the near-linear space regime by presenting a data structure that can be built in $\tilde{O}(n)$ time and answers HIA queries in $O(\log n / \log \log n)$ time. As a direct corollary, we obtain an $\tilde{O}(n)$-size data structure that maintains the LCS of a static string and a dynamic string, both of length at most $n$, in time optimal for this space regime.

The main ingredients of our approach are fractional cascading and the utilization of an $O(\log n / \log \log n)$-depth tree decomposition. The latter allows us to break through the $\Omega(\log n)$ barrier faced by previous works, due to the depth of the considered heavy-path decompositions.

1 Introduction

The solutions to algorithmic problems on texts frequently involve the construction of text indexes that can be built efficiently and offer a broad functionality, without significantly increasing space usage. A prime example of such an index is the suffix tree, which is ubiquitous in stringology. The work of Weiner [22] that introduced it, showed that it can be used to efficiently solve a number of fundamental open problems such as the computation of occurrences of patterns (given in an online manner) in a text or the computation of the longest common substring of two strings. However, it is usually the case that a suffix tree needs to first be augmented with other data structures before it can efficiently answer

1 The $\tilde{O}(\cdot)$ notation hides factors polylogarithmic in $n$. 

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more sophisticated queries, e.g., returning the longest common prefix of two substrings or the longest palindrome centered at some position; an augmentation with a lowest common ancestors data structure suffices for these examples [16,17].

Crucially, a text index, such as the suffix tree, is built once and can then be queried an arbitrary number of times. This is increasingly relevant: in many real-world scenarios, large pieces of information are stored on servers and are constantly queried by a large number of remote clients. From this perspective, it makes sense to devote some time to preprocess the data stored on the server in order to be able to provide quick responses to remote users later.

The Heaviest Induced Ancestors problem, which was introduced by Gagie et al. [14] and is defined next, has been proved to be useful in solving several variants of the problem of computing a longest common substring of two strings [1,4,5,8,14].

We say that a tree is weighted if there is a weight associated with each node $u$ of the tree, such that weights along root-to-leaf paths are increasing, i.e., for any node $u$ other than the root the weight of $u$ is larger than the weight of $u$’s parent. Further, we say that a tree is labelled if each of its leaves is given a distinct label from $[n]$, where $n$ is the number of leaves. As an example of a rooted, weighted, and labelled tree, consider the suffix tree of a string $S$, where $\$ does not occur in $S$, with the label of each leaf being the starting position of the corresponding suffix and the weight of each node being the length of the string it represents.

Definition 1. For two rooted and weighted trees $T_1$ and $T_2$ on $n$ leaves, we say that two nodes $u \in T_1$ and $v \in T_2$, are induced (by label $\ell$) if and only if there are leaves $x$ and $y$ labelled with $\ell$, such that $x$ and $y$ are weak descendants of $u$ and $v$, respectively.

Previous results and our contribution. Table 1 shows the state-of-the-art size vs. query-time tradeoffs for the HIA problem prior to our work and our result. Gagie et al. [14] presented several tradeoffs which have been since improved. We stress that the $\mathcal{O}(n \log^2 n)$-size data structure with query-time $\mathcal{O}(\log n)$ included in Table 1 was only sketched in [14]. We briefly discuss this sketch in Appendix A, as some of the ideas involved are similar to the ones we use. The remaining $\mathcal{O}(n)$-size known data structures found in Table 1 are due to Abedin et al. [1]. Charalamopoulos et al. [8] showed an unconditional lower bound for near-linear size data structures and a data structure with query-time $\mathcal{O}(1)$ and size $\mathcal{O}(n^{1+\epsilon})$ for any constant $\epsilon > 0$. We now formally state our main result, which matches the lower bound of [8].

Theorem 2. For any $\epsilon > 0$, there is an $\mathcal{O}(n \log^{2+2\epsilon} n)$-size data structure for the HIA problem that can be constructed in $\tilde{\mathcal{O}}(n)$ time and answers queries in $\mathcal{O}(\log n / \log \log n)$ time.

Applications of HIA. Before discussing some concrete applications of the HIA problem in string algorithms and the consequences of our results for them, we give a high-level description of how the HIA problem comes up in variants of computing an LCS.

Consider a string $S$ and a chosen subset $A$ of its positions, that we call anchors. Further, consider the following two tries: a trie $T^+$ for the strings in $\{S[1..k-1]^R : k \in A\}$, where $U^R$ denotes the reversal of $U$, and a trie $T^-$ for the strings in $\{S[k..|S|] : k \in A\}$. In other words, for every anchor $k \in A$, we have a path in the first trie for every prefix of $S[1..k-1]^R$ and a path in the second trie for every prefix of $S[k..|S|]$. We label each leaf of the two tries
Table 1 Size vs. query-time tradeoffs for the HIA problem; the size is measured in machine words.

<table>
<thead>
<tr>
<th>Size</th>
<th>Query time</th>
<th>Paper</th>
</tr>
</thead>
<tbody>
<tr>
<td>$\tilde{O}(n)$</td>
<td>$\Omega(\log n \log \log n)$</td>
<td>[8]</td>
</tr>
<tr>
<td>$O(n)$</td>
<td>$O(\log^2 n / \log \log n)$</td>
<td>[1]</td>
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<tr>
<td>$O(n \log n)$</td>
<td>$O(\log n \log \log n)$</td>
<td>[1]</td>
</tr>
<tr>
<td>$O(n \log^2 n)$</td>
<td>$O(\log n)$</td>
<td>sketched in [14], see Appendix A</td>
</tr>
<tr>
<td>$O(n \log^{2+\epsilon} n)$</td>
<td>$O(\log n / \log \log n)$</td>
<td>this work</td>
</tr>
<tr>
<td>$O(n^{1+\epsilon})$</td>
<td>$O(1)$</td>
<td>[8]</td>
</tr>
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</table>

with the anchor it corresponds to. Now, observe that a substring $S[i..j]$ that crosses an anchor $k$, i.e., $i < k \leq j$, corresponds to an induced pair of nodes in the tries. Indeed, there is a path representing $S[i..k-1]^R$ in the first trie and a path representing $S[k..j]$ in the second trie. An illustration of this idea is provided in Figure 1. The set of anchors and the HIA queries performed in an application of this technique depends on the specific problem it is used for. For some of the usages, one may consider using a compressed form of tries [18].

As a first application, consider the maintenance of an LCS of a static string $T$ and a dynamic string $S$. By plugging our HIA data structure into the approach of [8], we obtain the following result, improving the state-of-the-art by polyloglog $n$ factors, and matching the lower bound for the update-time when nearly-linear space is available [8, Theorem 1].

▶ Corollary 3. We can maintain an LCS of a dynamic string $S$ and a static string $T$, each of length at most $n$, in $O(\log n / \log \log n)$ time per substitution operation using $\tilde{O}(n)$ space, after an $\tilde{O}(n)$-time preprocessing.

Further, the authors of [14] (implicitly) reduced to the HIA problem, the problem of preprocessing a text given in LZ77 compressed form so that one can compute its LCS with uncompressed patterns given online. Our HIA data structure yields the following result.

▶ Corollary 4. Let $S$ be a string of length $N$ whose LZ77 parse consists of $n$ phrases. We can store $S$ in $O(n \log N + n \text{polylog} n)$ space such that, given a pattern $P$ of length $m$, we can compute the LCS of $S$ and $P$ in $O(m \log n / \log \log n)$ time. For each pattern $P$, the returned result may be (consistently) incorrect with probability inverse polynomial in $n$.

Randomization is only used in the construction; all queries for the same pattern give identical results.

Figure 1 An illustration of the anchoring technique for LCS computation, with the constructed tries $T^-$ and $T^+$ drawn so that their roots are attached (in the middle). Any substring anchored at $k$, can be obtained by reading in a left-to-right manner the edge-labels from some node $u \in T^-$ to some node $v \in T^+$, that both have a leaf-descendant labelled with $k$.
Other applications of the HIA problem in string algorithms can be found in [1].

**Tree Decompositions.** One of the obvious divide-and-conquer techniques for efficiently solving algorithmic problems on trees is that of decomposing the tree(s) into smaller pieces and treating each of them separately. The most important attributes of a tree decomposition are usually its depth, i.e., the maximum number of pieces that one path can intersect, and the structure of each individual piece (e.g., pieces being paths may offer an advantage). We next describe some tree decompositions for a tree $T$ with $n$ nodes. For a node $v$, denote by $s(v)$ the number of nodes in $v$’s subtree.

Arguably, the most well-known tree decomposition is the heavy-path decomposition [17]. Abstractly, this decomposition is a partition of the edges into light and heavy, such that:

- all connected components after deleting the light edges are paths, called heavy paths;
- each root-to-leaf path consists of $O(\log n)$ prefixes of heavy paths and $O(\log n)$ light edges, i.e., the depth of the decomposition is $O(\log n)$.

A heavy-path decomposition can be realized in several ways; two of which are as follows:

- **HP1:** Each non-leaf node $u$ of the tree chooses a child $v$ with maximum $s(v)$ and the edge from $u$ to $v$ is designated as heavy. The remaining edges outgoing from $u$ are light.
- **HP2:** An edge $(u, v)$ is designated as heavy if and only if $\lfloor \log s(u) \rfloor = \lfloor \log s(v) \rfloor$.\(^3\)

Intuitively, using a heavy-path decomposition, one may often lift an algorithm that only works for paths and/or balanced trees to work for arbitrary trees – usually with some overhead.

All previous works on the HIA problem used heavy-path decompositions, which, as discussed, are of depth $\Omega(\log n)$. This adversely affects their query times as one may have to traverse the decomposition along a root-to-leaf path at query time. Thus, in order to achieve sublogarithmic query time, we considered tree decompositions of smaller depths. There are a couple of generalizations of heavy-path decompositions that have the sought depth, i.e., $O(\log n/\log \log n)$. We next discuss two such decompositions that are also based on partitioning the edges into light and heavy. The caveat is that, for each of them, the connected components after the removal of the light edges are trees, which we call heavy trees, instead of paths and hence some extra work is required.\(^4\)

The heavy $\alpha$-tree decomposition, introduced by Bille et al. [6], is of depth $O(\log_{\alpha} n)$ and is defined analogously to HP1: each non-leaf node $u$ chooses its (at most) $\alpha$ heaviest (with respect to subtree-sizes) children; the edge from $u$ to each of these children is designated as heavy, while all remaining edges outgoing from $u$ are designated as light. By setting $\alpha = \lfloor \log n \rfloor$ one gets the sought depth.

An alternative is the so-called ART decomposition due to Alstrup et al. [2], which, for an input integer parameter $b$, has depth $O(\log_b n)$. For ease of presentation, we consider $b$ to be equal to $\lfloor \log n \rfloor$ so that the depth of the decomposition is $O(\log n/\log \log n)$. A partition of the edges yields such an ART decomposition if and only if each heavy tree contains $O(\log n)$ nodes that have more than one child (in the heavy tree). Alstrup et al. [2] showed how to compute an ART decomposition by computing a set of leafmost light edges (in the spirit of

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3 In some works this has been called a centroid decomposition [10]. It should not be confused with the hierarchical decomposition of the tree obtained by recursively deleting a centroid node, that is, a node whose removal splits the tree into three roughly equal components [7, 15].

4 Heavy trees are sometimes called micro trees, while the tree obtained from $T$ by contracting each micro tree is called a macro tree. We avoid this notation to not confuse with the so-called micro-macro decomposition [5], which, for a positive integer $k \leq n$, is a partition of the vertices of $T$ into $O(n/k)$ sets, such that each set $S$ is of size $O(k)$ it induces a subtree of $T$ and has at most two vertices that have neighbours that are not in $S$. 
HP2 with the base of the logarithm changed from 2 to $\lceil \log n \rceil$, removing them along with their descendants from the tree, and recursing. Here, for convenience, we compute an ART decomposition similar to the HP2-realization of a heavy-path decomposition: an edge $(u, v)$ is heavy if and only if both $s(u)$ and $s(v)$ are in $(n/\lfloor \log n \rfloor^{k+1}, n/\lfloor \log n \rfloor^k]$. For each heavy tree, we call branches its maximal down-the-tree paths in which all nodes except the deepest one have exactly one child (in the heavy tree); each heavy tree has $O(\log n)$ branches.

Our techniques. In order to answer an HIA query for nodes $u$ and $v$, we consider $O(\log n/\log \log n)$ pairs of heavy trees that consist of a heavy tree in the root-to-$u$ path and a heavy tree in the root-to-$v$ path. For each such pair, we compute an induced pair $(x, y)$ of nodes in these heavy trees that are ancestors of $u$ and $v$, respectively, and have maximum total weight. Similarly to previous work, we observe that not every pair of heavy trees needs to be considered. Instead, it suffices to consider a number of pairs of heavy trees linear to the depth of the tree decompositions by a procedure analogous to the natural algorithm for checking whether there are two elements of a sorted list that sum to a target $t$: start with two pointers, one at the beginning of the list and one at the end and move each of them in only one direction (either to the right or left). For each pair of heavy trees, we construct data structures that can efficiently handle each of the cases of how the locations of the lowest ancestors of $u$ and $v$ in the heavy trees relate to the locations of the lowest ancestors (in the heavy trees) of same-label leaves. Each data structure considers similar cases as previous work, however now we are working with two trees instead of two paths, and hence need to be more careful. This way, we reduce an HIA query to $O(\log n/\log \log n)$ predecessor queries. By answering each of these predecessor queries independently, we obtain a data structure that answers HIA queries in $O(\log n)$ time. Indeed, in our case, each predecessor query requires $\Omega(\log \log n)$ time to be answered independently.

However, crucially, we show how to design the data structures so that all predecessor queries need only two values: the preorder number of $u$ or the preorder number of $v$. This is achieved by reordering the trees so that heavy edges come last. Then, larger preorder numbers correspond to a larger depth of the lowest ancestor on a branch. This means that the combination of our techniques with fractional cascading would yield a faster algorithm for answering all the predecessor queries; the first one for each queried value would take $O(\log \log n)$ time, while all subsequent ones would take $O(1)$ time each. The final technical hurdle is that fractional cascading requires the so-called underlying catalog graph to have polylogarithmic degree [21]. The construction of such a graph is straightforward if $T_1$ and $T_2$ are of polylogarithmic degree: roughly speaking, it suffices to consider the Cartesian product of two trees whose nodes represent branches and heavy trees of each of $T_1$ and $T_2$. We overcome this difficulty in the general case by reducing the maximum degree of these trees prior to taking their Cartesian product while maintaining all of their desirable properties.

2 Preliminaries

We use $[n]$ to denote the set $\{1, 2, \ldots, n\}$. Throughout the paper, we perform the same operations on $T_1$ and $T_2$ and define objects in these trees, so we are going to use $T_i$ to denote any of the trees. Similarly, we are going to use $v_*$ to denote a node $v_1 \in T_1$ or a node $v_2$ in $T_2$ etc. as an abbreviation of writing that some property holds for $v_i$ for both $i \in \{1, 2\}$.

Lowest Common Ancestor. LCA queries can be answered in constant time after an $O(n)$-time preprocessing [17].
8.6 Optimal Near-Linear Space Heaviest Induced Ancestors

**LowestCommonAncestor (LCA)**

**Input:** A rooted tree $T$.

**Query:** What is the node of largest depth that is an ancestor of both $u$ and $v$?

**Predecessor query.** For a static set $S$, a combination of $x$-fast tries [23] and deterministic dictionaries [20] yields an $O(n)$-size data structure that can be built in $O(n)$ time and answers predecessor queries in $O(\log \log U)$ time deterministically (cf., [12, Proposition 2]); this is optimal [19].

**PredecessorQuery**

**Input:** A set $S$ of $n$ integers from $[U]$.

**Query:** For a given integer $y$, what is the largest $x \in S$ such that $x \leq y$?

**Range Minimum Query.** RMQs can be answered in constant time after an $O(n)$-time preprocessing [11, 13]. By setting, for each $i$, $S[i] := |U| - S[i]$, we get a structure for the symmetric RangeMaximumQuery problem.

**RangeMinimumQuery (RMQ)**

**Input:** A sequence $S$ of $n$ integers from $[U]$.

**Query:** For given positions $i$ and $j$, with $1 \leq i \leq j \leq n$, what is (the position of) the minimum among $S[i], S[i + 1], \ldots, S[j]$?

**Deterministic Static Dictionary.** A dictionary is a structure that stores a set of keys (often with associated values) and allows answering membership queries (or getting the value of a given key). There are multiple randomized solutions, but there is even a deterministic solution with $O(n)$ space, $\tilde{O}(n)$-time preprocessing, and constant-time queries [20].

**Fractional cascading.** Consider a directed graph $C$, called the catalog graph, which has a sorted list (also called a catalog) in each of its nodes. Let the total size of the lists be $n$. Now, suppose that we want to answer queries of the following type: for a connected subgraph $G$ of $C$ and a query value $v$, find the predecessor of $v$ in each of the lists stored in the nodes of $G$.

A naive way of solving this problem would be to ignore any preprocessing and run a separate binary search in the sorted list of each of the nodes of $G$, for a total of $O(|G| \log n)$ time. Fractional cascading is a general optimization technique that allows the speed-up of multiple binary searches for the same value over multiple related sorted sequences of objects.

If the degree of each node of the catalog graph is bounded by a constant, the original solution of Chazelle and Guibas [9] answers a query in $O(\log n + |G|)$ time after a linear-time preprocessing in the comparison model. To be precise, it is sufficient for the catalog graph to have locally bounded degree (as per Definition 1 of [9]). Unfortunately, in our case, this is not useful. A subsequent work of Shi and JáJá [21] achieved the same complexities for graphs of polylogarithmic maximum degree in the word RAM model of computation (in the original description, the graph is a tree, however, there is no difficulty in extending this to the general setting considered by Chazelle and Guibas [9]). Crucially, these data structures can also handle the case where the nodes of $G$ are given one by one in an online manner; the only requirement is that each node (other than the first) must be a neighbor of some previous one. Note that the $O(\log n)$ term in the complexities comes from performing a binary search in the first of the considered lists. Then, the predecessor of the query value $v$ in each of the subsequently considered lists is obtained by following a constant number of
pointers, which can be retrieved in $O(1)$ time. (During the preprocessing phase, the catalogs are augmented in an appropriate manner and said pointers are constructed.) In the word RAM model of computation, the first query can be solved faster using other data structures, e.g., in $O(\log \log U)$ time with the structure discussed above for the PredecessorQuery problem.

3 An $\tilde{O}(n)$-size Data Structure with Optimal Query Time

Let $b > 1$ be a parameter to be chosen later. Consider a rooted, weighted, and labelled tree $T$. The weight of a node $u$ is denoted by $\text{weight}(u)$. For a node $v$, we denote by $s(v)$ the number of nodes in $v$’s subtree, including $v$. For an integer $k$, a node $v$ is on layer $k$ if and only if $n/b^k < s(v) \leq n/b^{k+1}$. An edge that connects nodes of the same layer is called heavy and the other edges are light. Each maximal subtree that does not contain any light edges is called a heavy tree. We stress that a heavy tree might be a singleton.

We decompose each heavy tree into branches, that is, maximal down-the-tree paths of nodes, where every node apart from the deepest one has one child (in the heavy tree). The last node is either a leaf of the heavy tree or has at least two children. Note that there can be branches consisting of a single node. We call a node implicit if it is an internal (non-leaf) node of the heavy tree with one child; otherwise, we call it explicit. For a heavy tree, we obtain a compacted version of it, called compacted heavy tree, by eliminating all the implicit nodes through the contraction of either of their incident edges. See an example in Figure 2.

Figure 2: An example tree with $n = 26$ is shown. Each node is labelled with its subtree size, while $b = 3$. On the left, the heavy edges are thick, while the light edges are thin. On the right, the tree is decomposed into heavy trees (marked with ellipses). The compaction of heavy trees is illustrated as follows: empty circles denote implicit nodes, while full circles denote explicit nodes.

Observe that in a compacted tree, every non-leaf node has at least two children. Hence, there are fewer internal nodes than there are leaves of the tree. As every leaf in a heavy tree has a sufficiently big subtree underneath it, we obtain a bound on the total number of nodes inside a single compacted heavy tree; this shows that we obtain an ART decomposition [2].

Lemma 5. There are at most $O(\log n / \log b)$ layers in a tree $T$ on $n$ leaves and each heavy tree has $O(b)$ branches.

Proof. Consider a heavy tree $H$ of $T$ and its compacted form $H_C$. Let $r$ be the root of $H$. For every leaf $\ell$ of $H$, we have that $s(\ell) > s(r)/b$ as $\ell$ and $r$ are nodes of the same heavy tree, and hence they are in the same layer of $T$. As the subtrees of $T$ rooted at the leaves of $H$ are
disjoint and their total size is at most \( s(r) \), \( H \) contains at most \( b \) leaves. Further, as there are no internal nodes with one child in \( H_C \), there are at most \( b - 1 \) non-leaf nodes in \( H_C \), and hence \( H_C \) has \( O(b) \) nodes. All branches in \( H \) are disjoint, so there are \( O(b) \) of them.

Consider the \( k \)-th layer of the tree \( T \) and a node \( u \) in this layer. As the subtree of \( u \) has at least one node, we have \( 1 \leq s(u) \leq n/b^k \), so \( k \leq \log_b n = \log n / \log b \).

For an HIA query \((v_1, v_2)\), the paths from \( v_1 \) and \( v_2 \) to the roots of their respective trees are called query paths. The result of an HIA query is a pair of nodes \((u_1, u_2)\), that are on the query paths from \( v_1 \) and \( v_2 \), respectively. To answer a query \((v_1, v_2)\), we identify the sequences of heavy trees \( B_1 \) and \( B_2 \) that contain nodes on the query paths from \( v_1 \) and \( v_2 \), respectively, and perform restricted HIA queries for some pairs of those heavy trees. More precisely, in each step of the algorithm, having chosen two heavy trees \( B_1[i] \) and \( B_2[j] \), we try to find the pair of induced ancestors \((u_1, u_2) \in B_1[i] \times B_2[j] \) of \((v_1, v_2)\) with the maximum combined weight or determine that there is no such pair. A pseudocode for this procedure is given as Algorithm 1.

Algorithm 1 Algorithm for answering HIA queries.

```
1 function hia(T_1, T_2, v_1, v_2)
2     (r_1, r_2) ← null
3     (B_1, B_2) ← sequences of heavy trees on paths from v_1 and v_2 down-the-tree
4     i ← 0
5     j ← |B_2| - 1
6     repeat
7         x_1 ← lowest ancestor of v_1 in B_1[i]
8         x_2 ← lowest ancestor of v_2 in B_2[j]
9         (u_1, u_2) ← restricted-hia(v_1, v_2, B_1[i], B_2[j], x_1, x_2)
10        if (r_1, r_2) = null or weight(u_1) + weight(u_2) > weight(r_1) + weight(r_2)
11            then
12                (r_1, r_2) ← (u_1, u_2)
13        if (i = |B_1| - 1) and (j = 0) then
14            return (r_1, r_2)
15        if i + 1 < |B_1| and roots of B_1[i + 1] and B_2[j] are induced then
16            i ← i + 1
17        else
18            j ← j - 1
19    until false
```

To make the description more modular, we provide \( x_1 \) and \( x_2 \) to the restricted HIA query, where \( x_1 \) is the lowest ancestor of \( v_1 \) in the considered heavy tree. Note that \( x_1 \) is either the parent of the root of \( B_1[i + 1] \) if \( i + 1 < |B_1| \) or \( x_1 \) equals to \( v_1 \) otherwise, and similarly for \( x_2 \). We note that in Algorithm 1, we ask restricted HIA queries about the same pair \((v_1, v_2)\) for various pairs \((B_1[i], B_2[j])\) of heavy trees and it may be that \( v_1 \notin B_1[i] \) and/or \( v_2 \notin B_2[j] \), whereas we also provide nodes \( x_1 \), in which \( v_1 \) connects to the respective heavy subtree in \( B_1 \).

Lemma 6. Algorithm 1 performs \( O(\log n / \log b) \) restricted HIA queries to find the heaviest induced ancestors of nodes \( u \) and \( v \).
Proof. The paths from \(v_i\) to the roots pass heavy trees with monotonically decreasing indices of layers, so the sequences \(B_1\) and \(B_2\) contain at most \(O(\log n / \log b)\) elements. After every restricted HIA query, we either increase \(i\) or decrease \(j\) by 1, so the total number of restricted HIA queries that we perform is at most \(|B_1| + |B_2| = O(\log n / \log b)\).

The correctness follows from the monotonicity of being induced. For an induced pair of nodes, any pair of their (weak) ancestors is also induced. Conversely, if a pair is not induced, any pair of their (weak) descendants is also not induced.

This implies that if the roots of \(B_1[i]\) and \(B_2[j]\) are induced then there is no need to query for any pair \((B_1[i'], B_2[j'])\) with \((i' < i) \land (j' < j)\), as such a query would return a pair of (strict) ancestors of the pair returned by the restricted HIA query for \((B_1[i], B_2[j])\). We call a pair of trees \((B_1[i'], B_2[j'])\) dominated, if there exist \(i > i'\) and \(j > j'\) such that the roots of \(B_1[i]\) and \(B_2[j]\) are induced. We show that the algorithm performs restricted HIA queries at Line 9 exactly for those pairs of heavy trees that (i) are not dominated and (ii) for which the result of the restricted HIA query is not null.

The algorithm maintains the invariant that each of the restricted HIA queries is called for the pairs of trees for which their roots are induced and that the pair \((B_1[i], B_2[j])\) is not dominated. This is true for the first iteration, where the pair \((B_1[0], B_2[|B_2| - 1])\) is considered, because \(B_2[|B_2| - 1]\) has a leaf and the root of \(B_1[0]\) is the root of \(T_1\). Now we show that the invariant is maintained later. We start with the assumption that the result of restricted-hia\((v_1, v_2, B_1[i], B_2[j], x_1, x_2)\) is not null and this pair is non-dominated. Now, we have to distinguish between two cases.

Case 1: We next consider pair \((B_1[i + 1], B_2[j])\). It means that the check in Line 15 confirmed that the roots of the the trees are induced, so the the result of calling restricted-hia\((v_1, v_2, B_1[i + 1], B_2[j], x_1, x_2)\) is not null. Further, this pair of heavy paths is not dominated since \((B_1[i], B_2[j])\) is not dominated.

Case 2: We next consider pair \((B_1[i], B_2[j - 1])\). This can only happen if \(i = |B_1| - 1\) or the roots of \(B_1[i + 1]\) and \(B_2[j]\) are not induced; in either of these cases, \((B_1[i], B_2[j - 1])\) is not dominated. Further, as the answer to the HIA query for pair \((B_1[i], B_2[j])\) is not null, the answer for \((B_1[i], B_2[j - 1])\) cannot be null either.

Thus, the invariant is maintained in both cases.

Clearly, the heaviest induced pair of ancestors of \(v_1, v_2\) belongs to a pair of heavy trees that satisfy conditions (i) and (ii). We claim that we process all such pairs. Observe that for a fixed \(j\) there are two indices \(0 \leq i_1 < i_2 \leq |B_1|\) such that the pair \((B_1[i_1], B_2[j])\) is dominated for \(0 \leq i < i_1\), non-dominated for \(i_1 \leq i < i_2\), and corresponds to a null answer for \(i_2 \leq i < |B_1|\). By the invariant, just after any decrease of \(j\) in Line 18 it holds that \(i \geq i_1\), as \((B_1[i], B_2[j])\) is non-dominated. Actually \(i = i_1\), because the pair \((B_1[i - 1], B_2[j])\) is dominated as in the previous step we considered the pair \((B_1[i], B_2[j - 1])\) for which the answer was not null. As in the next steps we process all \(i\) up to (but excluding) \(i_2\), the claim follows.

3.1 Restricted HIA Queries

In this subsection, we present a data structure that efficiently answers restricted HIA queries.

\begin{itemize}
\item \textbf{Theorem 7.} For every two trees \(T_1, T_2\) on \(n\) leaves and an integer parameter \(b \in \mathbb{N}\), there exists an \(O(n b^2 \log^2 n / \log^2 b)\)-size data structure that can be computed in \(\tilde{O}(n b^2 / \log^2 b)\) time and answers (i) queries about whether the roots of two given heavy trees are induced in constant time, (ii) any restricted HIA query \text{restricted-hia}(v_1, v_2, H_1, H_2, x_1, x_2)\) in
\end{itemize}
constant time plus the time required to answer a predecessor query about \( \text{pre}(v_1) \) and one about \( \text{pre}(v_2) \); these predecessor queries are performed on two out of \( \mathcal{O}(b^2) \) (preprocessed) lists stored for the pair of heavy trees \((H_1, H_2)\).

We divide the proof into three parts: we first describe the preprocessing phase, then discuss the properties of the created data structure, and, finally, present the query procedure.

**Preprocessing.** First, we compute the partition of the edges of each of \( T_1, T_2 \) into heavy and light, and the implied heavy trees. For each node, we store its assignment to the heavy tree and to the branch to which it belongs. Further, for each \( T_u \), we build a linear-size data structure for answering LCA queries in \( \mathcal{O}(1) \) time [17]. For each node in \( T_u \), we fix the order of its children such that the children that are in the same heavy tree are last. (The order of children that are connected to the parent with the same type of edge, heavy or light, is arbitrary.) Next, we compute preorder traversals of \( T_u \), for each node \( u \), we denote by \( \text{pre}(u) \) the preorder number of \( u \) and by \( T_u[p] \) the node of \( T_u \) whose preorder number is \( p \); we have \( T_u[\text{pre}(u)] = u \). Additionally, for each label, we identify the leaves of \( T_1 \) and \( T_2 \) with that label (recall the labels in a single tree are pairwise distinct).

Next, for each pair \((\ell_1, \ell_2)\) of leaves with the same label, we iterate over all pairs \((B_1^{\ell_1}[i_1], B_2^{\ell_2}[i_2])\) of heavy trees on their query paths and insert a point to the data structure for each pair of branches in \( B_1^{\ell_1}[i_1] \times B_2^{\ell_2}[i_2] \). This procedure is formalized as Algorithm 2.

**Algorithm 2** Preprocessing for a pair of leaves \((\ell_1, \ell_2)\) with the same label.

```plaintext
1 procedure add_label(T_1, T_2, \ell_1, \ell_2)
2     \((B_1^{\ell_1}, B_2^{\ell_2}) \leftarrow \text{sequence of heavy trees on query paths from } \ell_1 \text{ and } \ell_2 \)
3     for \( i_1 \leftarrow 0, 1, \ldots, |B_1^{\ell_1}| - 1 \) do
4         for \( i_2 \leftarrow 0, 1, \ldots, |B_2^{\ell_2}| - 1 \) do
5             for \( e_1 \leftarrow \text{branch of } B_1^{\ell_1}[i_1] \) do
6                 for \( e_2 \leftarrow \text{branch of } B_2^{\ell_2}[i_2] \) do
7                     \( w_1 \leftarrow \text{LCA}(\ell_1, \text{lowest node on } e_1) \)
8                     \( w_2 \leftarrow \text{LCA}(\ell_2, \text{lowest node on } e_2) \)
9                     insert point \((\text{pre}(w_1), \text{pre}(w_2))\) to \( D_{B_1^{\ell_1}[i_1], B_2^{\ell_2}[i_2]}[e_1, e_2] \)
```

We call pairs \((B_1, B_2)\) of heavy trees that are processed by Algorithm 2 *relevant*. We first run this algorithm once just to record all relevant pairs of heavy trees, without inserting any points to any structures. We then sort the relevant pairs, remove duplicates, and construct a deterministic dictionary over them [20]. This allows us to check in constant time if the roots of two trees are induced because this is equivalent to checking if the pair of trees is relevant. For each relevant pair of heavy trees, we initialize an array \( D_{B_1, B_2} \) indexed by pairs of branches \((e_1, e_2)\), where \( e_1 \) is a branch in \( B_1 \) and \( e_2 \) is a branch in \( B_2 \). In each entry of the array, we create (store a pointer to) a data structure for the corresponding pair of branches. We then re-run Algorithm 2, inserting the points to the structures as needed, with the help of the deterministic dictionary built for relevant pairs of heavy trees. As each branch \( e \) belongs to a unique heavy tree, we often drop the superscript and write \( D[e_1, e_2] \) instead of \( D_{B_1, B_2}[e_1, e_2] \). We call a pair \((e_1, e_2)\) of branches *relevant* if and only if pair of their assigned heavy trees is relevant. By Lemma 5, every query path is decomposed into \( \mathcal{O}(\log n/\log b) \) parts on different layers and each heavy tree has \( \mathcal{O}(b) \) branches. Hence, for every pair \((\ell_1, \ell_2)\) of leaves with the same label, we insert a point to \( \mathcal{O}(b^2 \log^2 n/\log^2 b) \) structures.
Finally, for each relevant pair of branches \( (e_1, e_2) \), we perform the following postprocessing of structure \( D[e_1, e_2] \):

- Remove all points \((x, y)\) for which there exists another point \((x', y')\) such that \(x \leq x', y \leq y'\) and \((x, y) \neq (x', y')\). This can be done in \( \tilde{O}(|D[e_1, e_2]|) \) time by sorting the points and processing them in the left-to-right order.
- Let \( D_x[e_1, e_2] \) and \( D_y[e_1, e_2] \) be the sets of \( x\)- and \( y\)-coordinates of the remaining points, respectively. We build a data structure for the \textsc{PredecessorQuery} problem for each of \( D_x[e_1, e_2] \) and \( D_y[e_1, e_2] \) separately.
- We build a data structure for the \textsc{RangeMaximumQuery} problem for the points remaining in \( D[e_1, e_2] \) sorted by \( x\)-coordinate, where the weight of a point \((x, y)\) is \( \text{weight}(T_1[x]) + \text{weight}(T_2[y]) \).

We call the above stage the postprocessing of \( D[e_1, e_2] \).

To summarize the whole preprocessing stage for trees \( T_1 \), for each of the \( n \) labels we add \( \mathcal{O}(b^2 \log^2 n/\log^2 b) \) points to structures \( D[\cdot, \cdot] \), for a total number of \( \mathcal{O}(nb^2 \log^2 n/\log^2 b) \) points. The postprocessing of all the structures \( D[\cdot, \cdot] \) takes nearly linear time in their size and hence the total running time is \( \tilde{O}(nb^2/\log^2 b) \). The structures for the \textsc{PredecessorQuery} and \textsc{RangeMaximumQuery} problems have size linear in the number of elements they are built over and hence the total space is \( \mathcal{O}(nb^2 \log^2 n/\log^2 b) \).

**Properties of structures \( D[e_1, e_2] \).** In this paragraph, we show some properties of the structures \( D[e_1, e_2] \) that are useful for answering restricted HIA queries efficiently.

**Property 8.** For every pair \((w_1, w_2)\) added to \( B^{e_1}[i_1]_i B^{e_2}[i_2][e_1, e_2] \), \( w_1 \) is either on \( e_* \) or on the path from the highest node of \( e_* \) to the root of \( B^{e_*}[i_*] \).

**Proof.** Recall that \( w_* \) is the lowest common ancestor of \( \ell_* \) and the lowest node \( q \) on \( e_* \). Observe that as \( B^{e_*}[i_*] \) is on the query path from \( \ell_* \), the root \( r \) of \( B^{e_*}[i_*] \) is an ancestor of both \( \ell_* \) and \( q \). Hence, \( w_* \) lies on the \( r \)-to-\( q \) path, which directly yields the statement. ▶

Note that after the first step of postprocessing, \( D[e_1, e_2] \) satisfies the following property:

**Property 9.** After the postprocessing, for every pair \((e_1, e_2)\) of branches, after sorting the points of \( D[e_1, e_2] \) increasingly by \( x\)-coordinate, the sequence of points is also sorted decreasingly by \( y\)-coordinate.

Informally, we can now consider a one-dimensional problem, with points forming a sequence that can be efficiently navigated both in \( x\) and \( y\)-coordinates via predecessor queries.

We next show how the computed data structures \( D[e_1, e_2] \) enable us to answer restricted HIA queries efficiently.

**Answering a restricted HIA query.** We are now ready to present how to answer a restricted HIA query for a pair \((v_1, v_2)\) of nodes and heavy trees \( B_1 \) and \( B_2 \) on the query paths from \( v_1 \) and \( v_2 \). Let \((r_1, r_2) = \text{restricted-hia}(v_1, v_2, B_1, B_2, x_1, x_2) \) be a pair of ancestors of \( v_1 \) and \( v_2 \) within the trees \( B_1 \) and \( B_2 \) that are induced and have the maximum total weight. Recall that \( x_* \) is the lowest weak ancestor of \( v_* \) that is in \( B_* \) and let \( e_* \) be the branch containing \( x_* \). Further, let \( \ell \) be the label inducing \((r_1, r_2)\) and let leaves \( \ell \) share this label.

First, we show that we can find an induced pair of ancestors of \( v_1 \) and \( v_2 \) with the maximum combined weight using the structure \( D[e_1, e_2] \) before postprocessing. Then, we show that after the postprocessing stage, we can still retrieve the correct answer but more efficiently, by performing predecessor queries for \( \text{pre}(x_1) \) and \( \text{pre}(x_2) \). Finally, we show that
we can call predecessor queries for pre(v1) and pre(v2) instead of pre(x1) and pre(x2). The last step is not important for the correctness or efficiency of a single restricted HIA query but improves the complexity of Algorithm 1. Indeed, as all predecessor queries are for one of pre(v1) or pre(v2), we can use fractional cascading. We explain this final component in detail in Section 3.2.

Recall that in Algorithm 2, we insert point (pre(w1), pre(w2)) to D[e1, e2], where w⋆ = LCA(l⋆, lowest node on e⋆). In the proof of Property 8, we mention that w⋆ always belongs to B⋆ as the root of B⋆ is an ancestor of both l⋆ and the lowest node on e⋆. There are two possible relative locations of w⋆ and x⋆ within a heavy tree:

- l is below x, when w⋆ is a (not necessarily proper) descendant of x;
- l is attached above x, when w⋆ is a proper ancestor of x.

There are four cases for the relative locations of l with respect to x1 and x2:

Case 1: l is attached above x1 and x2,
Case 2: l is attached above x1 and l is below x2,
Case 3: l is attached above x2 and l is below x1,
Case 4: l is below x1 and x2.

We next treat each of these cases. For each of them, we retrieve the pair of induced ancestors of x1 and x2 with the largest total weight among all pairs of ancestors induced by a label l appropriately located with respect to x1 and x2. Each of these variants gives us a candidate pair for the restricted heaviest induced ancestors of x1 and x2. In the end, we return the candidate with the largest total weight. Similar case analysis was performed in previous solutions for the HIA problem, e.g., in [14].

**Lemma 10.** The answer to restricted-hia (v1, v2, B1, B2, x1, x2) can be retrieved from the information stored in D[e1, e2] before the postprocessing.

**Proof.** By Property 8, for every two points (pre(w1), pre(w2)) and (pre(w′1), pre(w′2)) added to D[e1, e2], we have that w1 is either a weak ancestor or a descendant of w′1, and similarly for w2 and w′2. Hence, the preorder numbers of the nodes correspond to their depths and we can check the ancestry relation by comparing them: for nodes u, u′ on a path, u is a weak ancestor of u′ if and only if pre(u) ≤ pre(u′).

Using this property, we show how to reduce each of the four cases listed above to finding a specific point in a particular rectangular subset of points. For now, we ignore the efficiency of the queries (a trivial implementation takes linear time) and focus on showing that the correct answer to the restricted HIA query can be retrieved from D[e1, e2] before the postprocessing.

Case 1: Every leaf l that is attached above x1 and x2 in nodes w1 and w2 makes the pair (w1, w2) a candidate result of the restricted HIA query. Hence we need to find a point (x, y) in D[e1, e2] such that x < pre(x1), y < pre(x2), and weight(T1[x]) + weight(T2[y]) is maximum. Then, (T1[x], T2[y]) is a restricted HIA candidate pair for (v1, v2).

Case 2: We need to find a point (x, y) ∈ D[e1, e2] such that x < pre(x1), y ≥ pre(x2) and weight(T1[x]) is maximized. Then, (T1[x], x2) is a restricted HIA candidate pair for (v1, v2).

Case 3: This case is symmetric to Case 2. We need to find a point (x, y) ∈ D[e1, e2] such that x ≥ pre(x1), y < pre(x2) and weight(T2[y]) is maximized. Then, (x1, T2[y]) is a restricted HIA candidate pair for (v1, v2).

Case 4: We need to check if there exists a point (x, y) such that x ≥ pre(x1) and y ≥ pre(x2).

If so, the pair (x1, x2) is a restricted HIA candidate pair for (v1, v2).
In each of the cases, we return a pair, if one exists, of induced ancestors of \((x_1, x_2)\) and hence also of \((v_1, v_2)\). The label \(\ell\) of leaves \(\ell_1\) and \(\ell_2\) that induces the pair \((r_1, r_2)\) of heaviest induced ancestors of \((v_1, v_2)\) in \(B_1 \times B_2\) inserted the point to \(D[e_1, e_2]\), since \(B_1\) and \(B_2\) are on the query paths from \(\ell_1\) and \(\ell_2\), respectively. Hence, the pair \((r_1, r_2)\) is found while considering one of the four cases.

Now, we show that it suffices to run the above algorithm only for the points in \(D[e_1, e_2]\) after the postprocessing stage.

\begin{lemma}
The answer to restricted-hia \((v_1, v_2, B_1, B_2, x_1, x_2)\) can be retrieved from the information stored in \(D[e_1, e_2]\) after the postprocessing.
\end{lemma}

\textbf{Proof.} As discussed in the proof of Lemma 10, for any two points \((\text{pre}(w_1), \text{pre}(w_2))\) and \((\text{pre}(w'_1), \text{pre}(w'_2))\) added to \(D[e_1, e_2]\), \(w_1\) and \(w'_1\) lie on a single root-to-leaf path, so their preorder numbers are in the same order as their depths in the tree. Recall that the trees are monotonically weighted, that is, the weights along each root-to-leaf path are increasing, so we have that \(\text{pre}(w_1) \leq \text{pre}(w'_1)\) implies \(\text{weight}(w_1) \leq \text{weight}(w'_1)\).

Let \(w = (\text{pre}(w_1), \text{pre}(w_2))\) be the point corresponding to the answer found by the algorithm presented in Lemma 10. Note that the returned induced pair of ancestors is not necessarily \((w_1, w_2)\), e.g., it can be \((w_1, x_2)\). Suppose that \(w\) was removed during the postprocessing phase. If so, it happened because there exists a point \(w' = (\text{pre}(w'_1), \text{pre}(w'_2))\) where \(\text{pre}(w_1) \leq \text{pre}(w'_1)\) and \(w' \neq w\). If \(w'\) is processed in a different case than \(w\), then the pair of ancestors corresponding to \(w'\) has a larger total weight than the one returned, yielding a contradiction. If \(w'\) is processed in the same case as \(w\), then the pair of ancestors corresponding to \(w'\) gives a pair of ancestors whose total weight is not smaller than that of the returned pair. Hence, the reduction presented in the proof of Lemma 10 still holds for the set \(D[e_1, e_2]\) after the postprocessing.

Next, we present how to implement each of the four cases in Lemma 10 efficiently using the fact that the points in \(D[e_1, e_2]\) have been postprocessed.

\begin{lemma}
The answer to restricted-hia \((v_1, v_2, B_1, B_2, x_1, x_2)\) can be retrieved from the information stored in \(D[e_1, e_2]\) after the postprocessing, with two predecessor queries: one for \(\text{pre}(x_1)\) and one for \(\text{pre}(x_2)\).
\end{lemma}

\textbf{Proof.} By computing the predecessor of \(\text{pre}(x_1)\) in \(D_x[e_1, e_2]\), we obtain intervals \(T^{x<\text{pre}(x_1)}_x\) of \(D_x[e_1, e_2]\). Similarly, intervals \(T^{y<\text{pre}(x_2)}_y\) and \(T^{y>\text{pre}(x_2)}_y\) of \(D_y[e_1, e_2]\) are obtained by computing the predecessor of \(\text{pre}(x_2)\) in \(D_y[e_1, e_2]\). Note that by Property 9, points from \(T^{y<\text{pre}(x_2)}_y\) (resp. \(T^{y>\text{pre}(x_2)}_y\)) of \(D_y[e_1, e_2]\) correspond to points from the interval of \(D_x[e_1, e_2]\) that we denote \(T^{y>\text{pre}(x_2)}_x\) \((T^{y<\text{pre}(x_2)}_x)\). Hence, we can translate each of the conditions on points in the cases of Lemma 10 to an intersection \(I_x^{\text{Case } i}\) of two intervals on \(D_x[e_1, e_2]\). This reduces each of the four cases to:

\begin{itemize}
  \item \textbf{Case 1:} Find the point with maximum weight \(\text{weight}(T_1[x]) + \text{weight}(T_2[y])\) in \(I_x^{\text{Case } 1}\) using an RMQ.
  \item \textbf{Case 2:} By the monotonicity of weights with respect to \(x\)-coordinates, the point with maximum weight \(\text{weight}(T_1[x])\) in \(I_x^{\text{Case } 2}\) is the rightmost element of \(I_x^{\text{Case } 2}\).
  \item \textbf{Case 3:} By the monotonicity of weights with respect to \(y\)-coordinates and Property 9, the point with maximum weight \(\text{weight}(T_2[y])\) in \(I_x^{\text{Case } 3}\) is the leftmost element of \(I_x^{\text{Case } 3}\).
  \item \textbf{Case 4:} It suffices to check if \(I_x^{\text{Case } 4}\) is non-empty.
\end{itemize}
Lemma 13. The answer to restricted-hia \((v_1, v_2, B_1, B_2, x_1, x_2)\) can be retrieved from the information stored in \(D[e_1, e_2]\) after the postprocessing, with two predecessor queries: one for \(\text{pre}(v_1)\) and one for \(\text{pre}(v_2)\).

Proof. Recall that in the approach presented in Lemma 12 we compute the predecessor of \(\text{pre}(x_1)\) in \(D_x[e_1, e_2]\) in order to divide \(D_x[e_1, e_2]\) into \(I_x^{<\text{pre}(x_1)} \) and \(I_x^{\geq\text{pre}(x_1)}\) and that all elements in \(D_x[e_1, e_2]\) are of the form \(\text{pre}(w_1)\) for a node \(w_1\) on the path from the lowest node on \(e_1\) to the root of \(B_1\).

We consider only \(v_1\) and \(x_1\) as the analysis for \(v_2\) and \(x_2\) is symmetric. We can focus on the case where \(v_1 \neq x_1\), as the other case is immediate. We clearly have that \(\text{pre}(v_1) \geq \text{pre}(x_1)\) as \(v_1\) is a descendant of \(x_1\). Recall that in the preprocessing stage, we reordered children of every node in such a way that children connected by a light edge are before those connected by a heavy edge, so if there is a child \(x_1'\) of \(x_1\) on \(e_1\) we have \(\text{pre}(x_1) \leq \text{pre}(v_1) < \text{pre}(x_1')\), as \(v_1\) is a descendant of a light child of \(x_1\) (by the definition of \(x_1\)). Hence, the predecessor of \(\text{pre}(v_1)\) is the same as the predecessor of \(\text{pre}(x_1)\) in \(D_x[e_1, e_2]\).

This concludes the proof of Theorem 7.

Finally, by setting the value of \(b\) to \(\lfloor \log^\epsilon n \rfloor\) for any constant \(\epsilon > 0\), we obtain a data structure using \(O(n \log^{2+2\epsilon} n / (\log \log n)^2) = O(1/\epsilon^2 \cdot n \log^{2+2\epsilon} n / (\log \log n)^2)\) space capable of answering restricted HIA queries in constant time plus the time required for answering two predecessor queries: one for \(\text{pre}(v_1)\) and one for \(\text{pre}(v_2)\). Algorithm 1 performs \(O(\log n / \log b)\) restricted HIA queries in order to answer an HIA query, so the total time required is \(O(1/\epsilon \cdot \log n)\). However, as all the predecessor queries ask about one of two values in different lists that are related to each other, we can make use of fractional cascading.

We omit the \(1/\epsilon\) factor in further sections and use the \(O_{\epsilon}(\cdot)\) notation instead to indicate a dependency on \(\epsilon\).

3.2 Fractional Cascading

For most of the cases described in the previous subsection, our structures are issuing predecessor queries. This is the only reason why the time complexity of an HIA query with our approach is not yet \(O_{\epsilon}(\log n / \log \log n)\). We will exploit the fact that all these queries look for the same target value (\(\text{pre}(v_1)\)) for structures built for \(T_1\) and \(\text{pre}(v_2)\) for structures for \(T_2\) but for different pairs of branches, which enables us to use fractional cascading.

We can think of creating two catalog graphs from \(T_1 \times T_2\) with nodes representing pairs \((e_1, e_2)\) of branches, storing the contents of \(D_x[e_1, e_2]\) in one catalog graph and those of \(D_y[e_1, e_2]\) in the other one. The execution of Algorithm 1 can be then seen as the traversal of a path in such a graph where, for a pair \((B_1[i], B_2[j])\) of heavy trees for which a restricted HIA query is performed by the algorithm, we query the catalogs of the nodes representing the pair of branches \((e_1, e_2)\) that contain the lowest weak ancestors of \((v_1, v_2)\) that are in \(B_1[i]\) and \(B_2[j]\), respectively. The problem with this direct approach is that it is not guaranteed that the degree of all vertices in each catalog graph is polylogarithmic: we might need to move from the node corresponding to two branches \((e_1, e_2)\) to any node corresponding to two branches \((e_1', e_2')\), where the heavy tree containing \(e_1'\) is attached to \(e_1\), and there could be even \(\Omega(n)\) such branches \(e_1'\).

We need to create catalog graphs in which the length of the considered path for each HIA query is \(O_{\epsilon}(\log n / \log \log n)\), while the degree of each node is \(O(\text{polylog } n)\) in order to be able to apply the result of Shi and JáJá [21]. This would ensure that all predecessor queries in \(D_x[\cdot, \cdot]\) and \(D_y[\cdot, \cdot]\) take constant time, apart from the first ones, which take \(O(\log \log n)\)
time using an $x$-fast trie. We describe how to build the appropriate catalog graphs below. As the shape of the graph for $D_1[\cdot,\cdot]$ and $D_2[\cdot,\cdot]$ is the same and only the contents of catalogs differ, we will only describe how to build one of them.

We preprocess $T_1$ and $T_2$ separately, first to compute trees $B(T_1)$ and then to build catalog graphs $C(T_1)$. From this, we build the catalog graph $C$ that can be seen as a Cartesian product of $C(T_1)$ and $C(T_2)$. More precisely, each node in $C$ is a pair $(v, w)$ for $v \in C(T_1)$ and $w \in C(T_2)$. For an edge between nodes $v_1$ and $v_2$ in $C(T_1)$, in the final catalog graph, we create edges between nodes $(v_1, w)$ and $(v_2, w)$ for each $w \in C(T_2)$. Analogically, for an edge between nodes $w_1$ and $w_2$ in $C(T_2)$, in the final catalog graph we create edges between nodes $(v, w_1)$ and $(v, w_2)$ for each $v \in C(T_1)$. This way, if the degrees of $C(T_1)$ and $C(T_2)$ are polylogarithmic, so is the degree of $C$.

We now explain how to build tree $B(T_\ast)$ from $T_\ast$. $B(T_\ast)$ contains nodes representing heavy trees (called heavy tree nodes) and nodes representing branches (called branch nodes):

- for each heavy tree $H$, we connect all branch nodes representing branches in $H$ as children of the heavy tree node representing $H$,

- for each heavy tree $H$, except the tree containing the root of $T_\ast$, we connect the heavy tree node representing $H$ as child of the branch node representing the branch containing the parent of the root of $H$.

**Proposition 14.** The depth of $B(T\ast)$ is $\mathcal{O}(\log n/\log \log n)$ and each heavy tree node has $\mathcal{O}(\log^2 n)$ children.

Recall that every heavy tree has $\mathcal{O}(b) = \mathcal{O}(\log n)$ branches, so every heavy tree node has $\mathcal{O}(\log^2 n)$ children. Any root-to-leaf path in $B(T\ast)$ alternates between heavy tree nodes and branch nodes. For any root-to-$v$ path $p$ in $T\ast$, there is a corresponding path $p'$ in $B(T_\ast)$ that visits the heavy tree nodes that correspond to the heavy trees that intersect $p$ and the branch nodes for which Algorithm 1 (when called for a pair of nodes containing $v$) could call predecessor queries for $D_2[\cdot,\cdot]$ and $D_1[\cdot,\cdot]$. For any $p$, $p'$ is of length $\mathcal{O}(\log n/\log \log n)$ and can be found in $\mathcal{O}(|p'|)$ time by following the path from the heavy tree containing $v$ to the root of $B(T\ast)$.

We now describe how to create a catalog graph $C(T\ast)$ from $B(T\ast)$. The construction is recursive and follows from the proof of Lemma 15 applied with $d = \mathcal{O}(\log n/\log \log n)$ and $b = \lfloor \log^2 n \rfloor$. The idea is to replace the structure of children of branch nodes having too many children with appropriate gadgets that roughly preserve the structure of the tree, do not increase the depth of the tree asymptotically and reduce the degree of each node to $\mathcal{O}(\text{polylog } n)$. Due to Proposition 14, we do not need to alter the structure of children for heavy tree nodes.

**Lemma 15.** For any depth-$d$ tree $B(T\ast)$ with $\mathcal{O}(n)$ nodes and parameter $b$, there is a tree $C(T\ast)$ satisfying all the following conditions:

- $C(T\ast)$ has $\mathcal{O}(n)$ nodes,

- all nodes of $C(T\ast)$ have degree $\mathcal{O}(b)$,

- $C(T\ast)$ has depth $d + \mathcal{O}(\log n/\log b)$,

- for each simple path $p$ in $B(T\ast)$, we can compute in $\mathcal{O}(|p'|)$ time a simple path $p'$ in $C(T\ast)$, such that $p$ is a subsequence of $p'$ and $|p'| \leq d + \mathcal{O}(\log n/\log b)$.

**Proof.** Consider a (branch) node $e$ of $B(T\ast)$ whose children, read left-to-right, are heavy tree nodes $h_1, h_2, \ldots, h_k$ for $k > b$. We replace this subgraph that contains $k + 1$ nodes with a gadget graph whose root is $e$ and whose set of nodes is a superset of $\{e\} \cup \{h_i : i \in [k]\}$.
For each node $u$, let $s(u)$ be the number of nodes in the subtree of $u$ in the considered tree. Let $s_0 = 0$ and, for $i \geq 1$, let $s_i$ be the prefix sum $s(h_1) + s(h_2) + \ldots + s(h_i)$. If there is an integer $\ell$ such that $s_{i-1} < \ell \cdot s(e)/b$ and $s_i \geq \ell \cdot s(e)/b$, we mark $h_i$. We call each set of consecutive unmarked nodes an interval. As $\ell \leq b$, there are $O(b)$ marked nodes and $O(b)$ intervals.

We create a gadget for $e$ (and, recursively, for some other nodes created in the construction, as described later) as follows:

- We attach as a child of $e$ every node that is either marked or is the only element of its interval.
- For each interval of more than one node, we create a new node $i_j$, called an interval node, attach it as a child of $e$, and attach all the nodes of the interval as children of $i_j$.

We recursively apply the same construction for any of the newly created interval nodes $i_1, i_2, \ldots, i_m$ whose degree is larger than $b$. See Figure 3 in Appendix B for an illustration.

From the construction, it follows that the degree of each node of $C(T_*)$ is $O(b)$ and that the size of $C(T_*)$ is $O(n)$, as all new nodes are of out-degree at least 2.

Let $u$ be a node in $T_*$. We now show that the depth for a node $e_u \in C(T_*)$ representing a branch containing $u$ is at most $d + O(\log n / \log b)$ by considering the edges above $e_u$ in $C(T_*)$. The edges can be of two types:

- Edges that are incident to at least one node that is not an interval node. By the depth of $B(T_*)$, we have at most $d$ such edges.
- Edges between interval nodes. For each such edge $(v, z)$, we have $s(v) \geq s(z) \cdot b$. Thus, similarly to the proof of Lemma 5, there are $O(\log n / \log b)$ such edges on the path from the root of $C(T_*)$ to $e_u$.

This concludes the proof of the bound on the depth of $C(T_*)$.

Each simple path $p$ in $B(T_*)$ naturally corresponds to a simple path $p'$ in $C(T_*)$. In particular, for each edge $(v, w)$ in $B(T_*)$, one can explicitly store a path $C(T_*)$ to which $(v, w)$ corresponds. The concatenation of all such paths for edges on $p$ yields a simple path $p'$ in $O(|p'|)$ time. As the depth of $C(T_*)$ is $d + O(\log n / \log b)$, the bound on $|p'|$ follows.

From $C(T_1)$ and $C(T_2)$ constructed as in Lemma 15, we create the catalog graph $C$ as described earlier. Only nodes that represent pairs of branches contain non-empty original catalogs. After the original catalogs are filled, we run the preprocessing of fractional cascading and appropriate augmented catalogs are created for all nodes in $C$ as described in [9, 21]. The $O(\log n / \log \log n)$ predecessor queries coming from restricted HIA queries performed in Algorithm 1 are naturally reduced to a constant number of queries to the $x$-fast tries and the traversal of a path of length $O(x / \log \log n)$ in $C$. This takes $O(x \log n / \log \log n)$ time in total and concludes the description of our data structure and the proof of Theorem 2.

References


As mentioned in the introduction, Gagie et al. \cite{14} sketched an $O(n \log^2 n)$-size data structure that answers HIA queries in $O(\log n)$ time, in the last paragraph of Subsection 2.1 of their work. They construct a data structure for each pair of heavy trees of $T_1$ and $T_2$ and reduce an HIA query for nodes $u$ and $v$ to a predecessor query in the data structure of each of $O(\log n)$ pairs of heavy trees. The idea for improving the (fully-described) $O(\log n \log \log n)$-time procedure for answering queries with a more efficient one is similar to ours and involves fractional cascading. They would need to build a catalog graph with nodes being pairs of heavy paths, reduce its degree, and make sure that the predecessor queries have the same target (by asking for the preorder numbers of $u$ and $v$). This is similar to what we describe in Section 3.2 for a different tree decomposition.

**B** Omitted Figure from Section 3.2

\begin{figure}[h]
\centering
\includegraphics[width=0.8\textwidth]{figure3}
\caption{On the left there is a branch $e$ of $T_*$, with outgoing edges to $h_1, h_2, \ldots, h_k$ in $B(T_*)$. On the right, there is a catalog graph gadget created for $e$, which is part of $C(T_*)$. Circles denote interval nodes and rectangles heavy tree nodes. Some of the intervals are recursively replaced with the gadget to decrease their degree. Heavy tree nodes have degree $O(b)$. In $C(T_*)$, $e$’s parent is the heavy tree to which it belongs, while the children of each $h_i$ are the branches in $h_i$.}
\end{figure}
From Bit-Parallelism to Quantum String Matching for Labelled Graphs

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Abstract

Many problems that can be solved in quadratic time have bit-parallel speed-ups with factor $w$, where $w$ is the computer word size. A classic example is computing the edit distance of two strings of length $n$, which can be solved in $O(n^2/w)$ time. In a reasonable classical model of computation, one can assume $w = \Theta(\log n)$, and obtaining significantly better speed-ups is unlikely in the light of conditional lower bounds obtained for such problems.

In this paper, we study the connection of bit-parallelism to quantum computation, aiming to see if a bit-parallel algorithm could be converted to a quantum algorithm with better than logarithmic speed-up. We focus on string matching in labeled graphs, the problem of finding an exact occurrence of a string as the label of a path in a graph. This problem admits a quadratic conditional lower bound under a very restricted class of graphs (Equi et al. ICALP 2019), stating that no algorithm in the classical model of computation can solve the problem in time $O(|P||E|^{1-\epsilon})$ or $O(|P|^{1+\epsilon}|E|)$. We show that a simple bit-parallel algorithm on such restricted family of graphs (level DAGs) can indeed be converted into a realistic quantum algorithm that attains subquadratic time complexity $O(|E|\sqrt{|P|})$.

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Keywords and phrases Bit-parallelism, quantum computation, string matching, level DAGs

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1 Introduction

Exact string matching problem is to decide if a pattern string $P$ appears as a substring of a text string $T$. In the classical models of computation, this problem can be solved in $O(|P| + |T|)$ time [10]. Different quantum algorithms for this basic problem have been developed [13,14,16], resulting into different solutions, the best of which finds a match in $O(\sqrt{|T|}(\log^2 |T| + \log |P|))$ time [13] with high probability. These assume the pattern and text are stored in quantum registers, requiring thus $O(|P| + |T|)$ qubits to function. Moreover, these approaches may rely on applying a linear number of quantum gates in parallel on different qubits. For example, Niroula and Nam [13] perform $O(\log(|T|))$ rounds of parallel swaps, executing $O(|T|)$ swaps in parallel per round.
In the classical models of computation, an analogy for these assumptions is to assume that the text has been preprocessed for subsequent queries. For example, one can build a Burrows-Wheeler transform-based index structure for the text in time $O(|T|)$ [4], assuming $T \in \{1, 2, \ldots, \sigma\}^*$, where $\sigma \leq |T|$. Then, one can query the pattern from the index in $O(|P| \log \log \sigma)$ time [4, Theorem 6.2]. In this light, quantum models can offer only limited benefit over the classical models for exact string matching.

Motivated by this difficulty in improving linear-time solvable problems using quantum approaches, let us consider problems known to be solved in quadratic time. For example, approximate string matching problem is such a problem: decide if a pattern string $P$ is within edit distance $k$ from a substring of a text string $T$, where edit distance is the number of single symbol insertions, deletions, and substitution needed to convert a string to another. This problem can be solved using bit-parallelism in $O(|P|/w(|T|))$ time [11], under the Random Access Memory (RAM) model with computer word size $w$. A reasonable assumption is that $w = \Theta(\log |T|)$, so that this model reflects the capacity of classical computers. Thus, when $|P| = |T| = n$, this bit-parallel algorithm for approximate string matching takes time at least $\Omega(n^{2-\epsilon})$ for all $\epsilon > 0$, as $\log n = o(n^\epsilon)$ for all $\epsilon > 0$. It is believed that this quadratic bound cannot be significantly improved, as there is a matching conditional lower bound saying that if approximate pattern matching could be solved in time $O(n^{2-\epsilon})$ with some $\epsilon > 0$, then the Orthogonal Vectors problem itself can be solved in subquadratic time [4, Theorem 6.2].

In this quest for breaking the quadratic barrier, we study another problem with a bit-parallel solution and a conditional lower bound. Consider exact pattern matching on a DAG: for every two nodes $u$ and $v$, there is an $OVH$ lower bound conditionally refuting an $O(|P||E|^{1-\epsilon})$ or $O(|P|^{1-\epsilon}|E|)$ time solution [7]. This conditional lower bound holds even if graph $G$ is a level DAG: for every two nodes $u$ and $v$, holds the property that every path from $u$ to $v$ has the same length. On DAGs, this string matching on labeled graphs (SMLG) problem can be solved in $O(|P|/w(|E|))$ time [15] in the bit-parallel model, so the status of this problem is identical to that of approximate pattern matching on strings. However, the simplicity of the bit-parallel solution for SMLG on level DAGs enables a connection to quantum computation. We consider a specific model of quantum computation, the Quantum Random Access Memory (QRAM) model [8], in which we have access to “quantum arrays”, and we assume that integer values like $|P|$, $|V|$ or $|E|$ fit into a (quantum) memory word. Under this model, we turn the bit-parallel solution into a quantum algorithm that solves SMLG on level DAGs with high probability in $O(|E|/\sqrt{|P|})$ time, breaking through the classical quadratic conditional lower bound.

Classical conditional lower bounds are not new to be broken by quantum computing. For example, the quadratic Orthogonal Vectors problem itself can be solved in subquadratic time (linear using QRAM) using quantum computing. This is not the only problem to have a better-than-quadratic solution in the quantum realm [17]. Nevertheless, to the best of our knowledge, we are the first to propose a subquadratic time algorithm for SMLG, even if restricted to a specific class of graphs. Moreover, the translation of a bit-parallel strategy to a quantum-parallel one is an original technique, and we are not aware of any other work utilising it.
An earlier work [6] provided a quantum algorithm solving SMLG in time $O(\sqrt{|V||E|}|P|)$. When the graph is non-sparse, that is $|V| = O(\sqrt{|E|})$, the time complexity becomes $O(|E|^3/2|P|)$, which is an improvement over classical algorithms. We offer a different kind of trade-off, limiting ourselves to a special class of graphs, but obtaining a better time complexity. We also note that, even if no subquadratic classical algorithm exists for non-sparse graphs, the existing classical reduction from OV [7] produces a sparse level DAG, for which our quantum algorithm runs in subquadratic time.

As mentioned above, in some previous works [13, 14, 16](and references in [16]) algorithms have been proposed to solve string matching in plain text in the QRAM model, under the assumption that a large number of quantum gates, possibly linear, can be applied in parallel when acting on different qubits. We find this assumption to be too restrictive, as even the classical RAM model does not adopt it, since in such a model of computation many operations would become trivial. Instead, our algorithm works without the need for such an assumption.

The paper is structured as follows. We revisit exact pattern matching and derive a simple brute-force quantum algorithm for SMLG, which we later improve on level DAGs. This improvement is based on extending the Shift-And algorithm [3], whose quantum version we extend for level DAGs.

In what follows, we assume the reader is familiar with the basic notions in quantum computing as covered in textbooks [12].

## 2 Preliminaries

An alphabet $\Sigma$ is a set of characters. Throughout the paper we assume $\Sigma$ is ordered, i.e., for each $a, b \in \Sigma$ we can decide if $a < b$. A sequence $P \in \Sigma^n$ is called a string and its length is denoted $n = |P|$. We denote integers $i, i + 1, \ldots, j$ as interval $[i,j]$ and represent a string $P$ as an array $P[0..n-1]$, where $P[i] \in \Sigma$ for $0 \leq i \leq n-1$, as in this work all indexes start from 0. String $P[i..j]$ is called a substring and string $P[0..i]$ a prefix of $P$. With bit-vectors discussed next, we use 0-based indexing.

Let $B$ be a $w$-bit integer interpreted as string $B[0..w-1]$ from alphabet $\{0,1\}$ such that $B = \sum_{i=0}^{w-1} B[i] \cdot 2^i$. We call $B$ a bit-vector. Given two bit-vectors $B$ and $C$, we define the following Boolean operations $A = B \land C$, $O = B \lor C$, and $N = \neg B$ as follows: $A[i] = 1$ iff $B[i] = C[i] = 1$, $O[i] = 1$ iff $B[i] = 1$ or $C[i] = 1$, and $N[i] = 1$ iff $B[i] = 0$. When bit-vector content is visualized, we list the most significant bit first, i.e., $B[w-1]B[w-2]\cdots B[0]$. With this in mind, we define the left-shifts $L = B \ll k$ and right-shifts $R = B \gg k$ as follows: $L[i+k] = B[i]$ and $R[i] = B[i+k]$. Here values out of the domain of the bit-vectors are assumed to be 0. Logarithms are assumed to be in base two: $\log n = \log_2 n$.

In directed labelled graph (DAG) $G = (V,E,\ell)$, $V$ is the set of nodes, $E$ is the sets of vertices, and $\ell : V \rightarrow \Sigma$ is a labelling function that assigns a character of the alphabet to each node. We assume the nodes to be indexed as $v_0,v_1,\ldots,v_{n-1}$ in topological order, where $n = |V|$. For $v_i \in V$, $\ell(v_i)$ is its label. Set of nodes $\text{in}(v_i) = \{j \mid (v_j,v_i) \in E\}$ contains the indexes of the in-neighbours of $v_i$, and $D_i = |\text{in}(v_i)|$ is the in-degree of $v_i$. If, for $0 \leq d \leq D_i - 1$, $v_k$ is the $d$-th in-neighbour of $v_i$ according to the topological indexing that we defined above, we express this fact using notation $k = \text{in}_i(d)$, where $\text{in}_i : [0,D_i-1] \rightarrow [0,n-1]$. 
In this work, we study the problem of string matching in labelled graphs, that consists in finding a match for a pattern string $P[0..m-1]$ in a labelled graphs $G$ over alphabet $\Sigma$, where $P$ has a match in $G$ if there is a path $v_1,\ldots,v_k$ such that $P = \ell(v_1)\cdots\ell(v_k)$ (we also say that $P$ occurs in $G$, and that $v_1,\ldots,v_k$ is an occurrence of $P$). Notice that if $|P| = 1$, a classic visit of the graph solves the problem in linear time, thus we always assume $|P| \geq 2$.

**Problem 1 (String Matching in Labeled Graphs (SMLG)).**

**Input:** A labeled graph $G = (V,E,L)$ and a pattern string $P$, both over an alphabet $\Sigma$.

**Output:** True if and only if there is at least one occurrence of $P$ in $G$.

### 3 Quantum Notation and Preliminaries

In quantum computing, data is represented in quantum bits (qubits), the quantum analogue to classical bits. A qubit can be in two states, denoted as $|0\rangle = \begin{pmatrix} 1 \\ 0 \end{pmatrix}$ and $|1\rangle = \begin{pmatrix} 0 \\ 1 \end{pmatrix}$ but, unlike a classical bit, it can also be a linear combination of the two states, a superposition: $|\psi\rangle = \alpha|0\rangle + \beta|1\rangle$. The complex values $\alpha$ and $\beta$ are called the amplitudes of $|\psi\rangle$. Measuring a qubit in superposition will result in either $|0\rangle$ or $|1\rangle$ with probabilities $|\alpha|^2$ and $|\beta|^2$, respectively.

Note that this notation can easily be generalised to integer states $|\psi\rangle = \sum_i \alpha_i |i\rangle$, as in the following example

$$\sum_{i=0}^{3} \alpha_i |i\rangle = \alpha_0|0\rangle + \alpha_1|1\rangle + \alpha_2|2\rangle + \alpha_3|3\rangle,$$

where $I$ is a quantum register of at least two qubits.

We assume to have a quantum random access memory (QRAM) able to use a quantum register as an index to access classical data. Let $m_0, m_1, \ldots, m_{n-1}$ be the data stored in QRAM $M$. Given quantum register $I$, the operation that reads data from $M$ into quantum register $Q$ initialized to $|0\rangle$ using $I$ as index is defined as follows [8]:

$$\sum_{i=0}^{n-1} \alpha_i |i\rangle \otimes |0\rangle \xrightarrow{\text{QRAM read}} \sum_{i=1}^{n} \alpha_i |i\rangle |0\rangle \oplus m_iQ = \sum_{i=1}^{n} \alpha_i |i\rangle |m_i\rangle.$$  

Notice that this is a unitary operation, and thus reading the same data into the same register twice will reset such a register to the value it had before performing the reading operation. In terms of time complexity, the execution of the read operation is proportional to the number of qubits in quantum register $I$. Under the Word-QRAM model with memory-word size $O(\log n)$ for inputs of size $n$, we can assume to be able to perform a QRAM read operation in $O(1)$, because $O(\log n)$ qubits are enough for register $I$ to index an input of size $n$. Indeed, this reflect the same assumption of the classical Word-RAM model, where operations on memory words are assumed to be constant.
A quantum computer, with access to QRAM, can solve the problem of finding an exact match for a pattern string \( P \) into a text string \( T \) in time \( O(|P| \sqrt{|T|}) \), with high probability. We explain a simple solution to this problem. Let \( |T| = n \) and \( |P| = m \), then \( T = t_0t_1\cdots t_{n-1} \) and \( P = p_0p_1\cdots p_{m-1} \) are two strings defined over a binary alphabet, that is \( t_i, p_j \in \{0, 1\} \) for \( 0 \leq i \leq n-1 \) and \( 0 \leq j \leq m-1 \). We use qubits \( C_T \) and \( C_P \) initialized to \( |0\rangle \) to track the current characters of \( T \) and \( P \), and we assume to have the text and the pattern stored in qubits in the following way:

\[
|0\rangle_{C_P} |0\rangle_{C_T} |t_0\rangle_{T_0} |t_1\rangle_{T_1} \cdots |t_{n-1}\rangle_{T_{n-1}} |p_0\rangle_{P_0} |p_1\rangle_{P_1} \cdots |p_{m-1}\rangle_{P_{m-1}}.
\]

We also use auxiliary qubits \( A_{-1}, A_0, A_1, \cdots A_{m-1} \), and quantum registers \( I, J \), and \( Q \), all three of log \( n \) qubits. We initialize \( A_{-1} \) and \( Q \) to \( |1\rangle \), while \( A_0, A_1, \cdots A_{m-1} \), \( I \), and \( J \) are all initialized to \( |0\rangle \). We prepare quantum register \( I \) in an equally balanced superposition spanning all the text positions, that is \( |0\rangle_I \rightarrow 1/\sqrt{n} \sum_{i=0}^{n-1} |i\rangle_I \), assuming \( n \) to be a power of 2, without loss of generality. If this is not the case, we generate a superposition as large as the first power of two greater than \( n \), then standard techniques can be applied to handle the additional substates, as explained in Appendix A.

Each individual state \( |i\rangle \) in the superposition represents a computation starting at position \( i \) in the text. In each of these computations, we scan \( T[i..i + m - 1] \) and try to match each character with \( P[0..m - 1] \), storing the intermediate results of such comparisons in registers \( A_0, A_1, \cdots, A_{m-1} \). More precisely, at iteration \( j \), \( 0 \leq j \leq m - 1 \), we compute a logical \( \text{xor} \) between \( t_{i+j} \) and \( p_j \) storing the result in \( C_P \) via a \( CX \) gate with control \( C_T \) and target \( C_P \). Then, we apply a \( X \) gate to \( C_P \), which now stores \(|\text{\text{xor}}(t_{i+j} \oplus p_j)\rangle_{C_P} = |t_{i+j} = p_j\rangle_{C_P} \). At this point, we apply a Toffoli gate with controls \( C_P \) and \( A_{j-1} \), storing the value in target qubit \( A_j \). We now reset \( C_T \) and \( C_P \) to \( |0\rangle \) by applying to them the same gates again, but in reverse order. As last step in iteration \( j \), we increase both \( I \) and \( J \) by 1 by performing transformation \( 1/\sqrt{n} \sum_{i=0}^{n-1} |1\rangle_I |j\rangle_J \rightarrow 1/\sqrt{n} \sum_{i=0}^{n-1} |1\rangle_I |i+1\rangle_J |j+1\rangle_J \) (this of course requires two separate addition operations), where the addition is intended to be modulo \( 2^n \). This allows us to read the next character of the pattern at the next iteration.

After the last iteration, we can run Grover’s operator \([9]\) where the marked items are represented by \(|a_{m-1}, i\rangle_{A_{m-1}} = |1\rangle\), and then measure register \( |I\rangle \) to locate the ending position of a match. Of course, we do not know the exact number of marked items, and we address this issue by guessing the number of items and rerunning the whole algorithm a constant number of times. We illustrate the entire procedure in Algorithm 1.

The algorithm is correct because, after each iteration of the for loop, we correctly keep track of the positions of the text that are active matches for the current prefix of the pattern.
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\[ \begin{array}{cccccc}
I & C_T & C_P & A_0 & A_1 \\
1 & 1 & |B|A & 0 & 0 & + \hline
2 & 1 & |B|A & 0 & 0 & + \\
3 & 1 & |A|A & 1 & 0 & + \\
4 & 1 & |A|A & 1 & 0 & + \\
5 & 1 & |B|A & 1 & 0 & + \\
6 & 1 & |B|A & 0 & 0 & + \\
7 & 1 & |A|A & 0 & 0 & + \\
8 & 1 & |B|A & 1 & 0 & + \\
\end{array} \]

Figure 1: An example of the evolution of the superposition after one iteration of Algorithm 1. The first arrow represents the application of a \( CX \) gate with control \( T_1 \) and target \( P_1 \), and the application of a \( X \) gate on \( P_1 \). The second arrow represents the application of a Toffoli gate with controls \( P_1 \) and \( A_0 \), and target \( A_1 \). Intuitively, in the first step we are checking that \( T[i + j] = P[j] \); in the second step we combine the result of this check with the contribution of the previous iteration(s).

At iteration \( j \), we assume by induction that register \( A_{j-1} \) stores \( |a_{j-1,i}\rangle_{A_{j-1}} = |1\rangle \) if and only if \( T[i\ldots i + j - 1] = P[0\ldots j - 1] \). Gates \( CX \) and \( X \) compute \( \neg(T[i + j] \oplus P[j]) \) storing it in \( C_P \). We then apply the Toffoli gate with controls \( C_P \) and \( A_{j-1} \), and target \( A_j \), obtaining superposition \( \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |i\rangle |a_{j-1,i} \wedge T[i + j] = P[j]\rangle_{A_j} \). Thus, \( |a_{j,i}\rangle_{A_j} = |a_{j-1,i} \wedge T[i + j] = P[j]\rangle_{A_j} \) if and only if \( T[i..i + j] = P[0..j] \).

As mentioned above, we have to be careful in running Grover’s search algorithm at the end of Algorithm 1. We defer these details to the full proof of Theorem 5 in Appendix C. For now, we assume that we are able to retrieve with arbitrarily high probability \( 1 - (7/8)^c \) a marked substate representing a match. Combining this with Lemma 1, we obtain the claimed result.

**Theorem 2.** Given a text string \( T \), pattern string \( P \) and integer \( c > 0 \), Algorithm 1 finds a match for \( P \) in \( T \) in time \( O(c(|P| \sqrt{|T|})) \). If there is no match, the algorithm returns a negative answer with probability \( p = 1 \). If there is at least one match, the algorithm returns the index of the last position of a match with probability \( c > 1 - (7/8)^c \).

**Proof.** For the correctness, consider Lemma 1 where \( j = m = |P| \), which is the number of times we run the for loop. In this case, \( |a_{|P|,i}\rangle_{A_{|P|}} = |1\rangle \) if and only if \( T[i..i+|P|-1] = P[0..|P|] \). Thus, measuring these substates yields a correct solution. The details of how to perform such a measurement respecting the time complexity and probability of success are deferred to the full proof of Theorem 5 in Appendix C.

5 String Matching in Labeled Graphs

5.1 Quantum Brute-Force Algorithm for SMLG

In SMLG we are given pattern string \( P \) with characters in alphabet \( \Sigma \) and a node-labeled graph \( G = (V,E) \), with labelling function \( \ell : V \to \Sigma \). We are asked to find a path (or, actually, a walk) \( \pi = v_1, v_2, \ldots, v_{|P|} \) in \( G \) such that \( \ell(v_1) \circ \ell(v_2) \circ \ldots \circ \ell(v_{|P|}) = P \), where \( \circ \) denotes string concatenation.

One could try to obtain a quantum algorithm for SMLG by generalizing the idea we presented for plain text. The idea would be to list all possible paths of length \(|P| \) in the graph, and then mark those ones that are actual matches for \( P \). Unfortunately, the superposition
An algorithm for solving exact string matching in plain text that, using QRAM, achieves $O(|P|^{1/2})$ time complexity. The details of how to handle Grover’s search at the end are given in Theorem 5, whose full proof is deferred to Appendix C.

**Algorithm 1**

**Input:** Text $T$ stored as $|t_0\rangle_T |t_1\rangle_T \cdots |t_n-1\rangle_T$, pattern string $P$ stored as $|p_0\rangle_P |p_1\rangle_P \cdots |p_m-1\rangle_P$, integer $c$

**Output:** A position of $T$ where a match for $P$ ends, if any

```
for $c$ times do
    Initialize quantum register $I$, $J$, $A_0$, $A_1\cdots A_{m-1}$ as $|0\rangle_I |0\rangle_J |0\rangle_{A_0} |0\rangle_{A_1} |0\rangle_{A_{m-1}}$;
    Initialize quantum register $A_{-1}$ and $Q$ as $|1\rangle_{A_{-1}}$ and $|1\rangle_Q$;
    // Apply $H^\otimes\log n$ to register $I$
    $|0\rangle_I \rightarrow \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |i\rangle_I$;
    for $m$ times do
        // Read $T[i]$ in $C_T$ and $P[j]$ in $C_P$ using registers $I$ and $J$ as indexes
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |i\rangle_I |j\rangle_J |0\rangle_{C_T} |0\rangle_{C_P} \rightarrow \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |i\rangle_I |j\rangle_J |t_i\rangle_{C_T} |p_j\rangle_{C_P} |C_P;$
        // Apply CX with control $C_T$ and target $C_P$
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |i\rangle_{C_T} |p_j\rangle_{C_P} \rightarrow \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |i\rangle_{C_T} |t_i\rangle_{C_P} |p_j\rangle_{C_P};$
        // Apply $X$ to $C_P$
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |t_i\rangle_{C_P} |p_j\rangle_{C_P} \rightarrow \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |-(t_i \oplus p_j)\rangle_{C_P} = \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |t_i = p_j\rangle_{C_P};$
        // Apply Toffoli with controls $C_P$ and $A_{j-1}$, and target $A_j$
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |t_i = p_j\rangle_{C_P} |a_{j-1}\rangle_{A_{j-1}} |0\rangle_{A_j} \rightarrow$
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |t_i = p_j\rangle_{C_P} |a_{j-1}\rangle_{A_{j-1}} |(t_i = p_j) \wedge a_{j-1}\rangle_{A_j};$
        // Reset $C_T$ and $C_P$ to $|0\rangle$ via uncomputation
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |-(t_i \oplus p_j)\rangle_{C_P} \rightarrow \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |t_i \oplus p_j\rangle_{C_P};$
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |t_i\rangle_{C_T} |t_i \oplus p_j\rangle_{C_P} \rightarrow \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |t_i\rangle_{C_T} |p_j\rangle_{C_P};$
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |i\rangle_I |j\rangle_J |t_i\rangle_{C_T} |p_j\rangle_{C_P} \rightarrow$
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |i\rangle_I |j\rangle_J |t_i \oplus t_i\rangle_{C_T} |p_j \oplus p_j\rangle_{C_P} = \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |i\rangle_I |j\rangle_J |0\rangle_{C_T} |0\rangle_{C_P};$
        // Increment indexes $I$ and $J$
        $\frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |1\rangle_Q |i\rangle_I |j\rangle_J \rightarrow \frac{1}{\sqrt{n}} \sum_{i=0}^{n-1} |1\rangle_Q |i \oplus 1\rangle_I |j + 1\rangle_J;$
        Apply gate $Z$ to qubit $R_{n-1}$, so that the sign of the amplitude is flipped if $|r_{n-1, j}\rangle_{R_{n-1}} = |1\rangle$;
        Choose $K \in [0,|P|]$ uniformly at random;
        Run Grover’s iterate operator the optimal number of times assuming to have $K$ solutions, with the oracle function being lines 5–14 of this algorithm;
        Measure $R_{n-1}$ into classical register $R_{cl}$;
        if $R_{cl} = 1$ then
        Measure $I$ into classical register $I_{cl}$ and return $I_{cl}$
```

return no
would be as large as there are paths of length $|P|$, and thus the overall time complexity would be $O(|P|\sqrt{|V||P|})$. Moreover, an adjacency matrix would be needed to check the existence of edges between nodes in constant time, yielding a space complexity of $O(|V|^2)$ qubits. We conclude that more involved techniques are needed.

### 5.2 The Classical Shift-And Algorithm

We first introduce the classical shift-and algorithm [3] for matching a pattern against a text and generalize it to work on graphs. Then, we show how the bit-vector data structure of that algorithm can be represented as a superposition of a logarithmic number of qubits. This approach allows us to achieve better performances than the brute force algorithm.

In the shift-and algorithm, we use bit vector $B$ of the same length of pattern $P$ to represent which of its prefixes are matching the text during the computation. Assuming integer-alphabet $\Sigma$, we also initialize bidimensional array $M$ of size $|P| \times |\Sigma|$ so that $M[j][c] = 1$ if and only if $P[j] = c$, and $M[j][c] = 0$ otherwise. The algorithm starts by initializing vector $B$ to zero and array $M$ as specified above. Then, we scan whole text $T$ performing the next four operations for each $T[i]$, $i \in [0, n-1]$, where $M[*][c]$ represents the $c$-th column of $M$:

1. $B \leftarrow B + 1$;
2. $B \leftarrow B \land M[*][T[i]]$;
3. if $B[m - 1] = 1$, return yes;

Operation 1 sets the least significant bit of $B$ to 1, which is needed to test $P[0]$ against $T[i]$. Operation 2 computes a bit-wise and between $B$ and the column of $M$ corresponding to character $T[i]$. Remember that $M[j][T[i]] = 1$ means $P[j] = T[i]$, thus this operation leaves each bit $B[j]$ set to 1 if and only if it was already set to 1 before this step and the the $j$-th character of the pattern matches the current character of the text. At this point, if bit $B[m - 1]$ is set to 1 we have found a match for $P$, and Operation 3 will return yes. For the other positions, if bit $B[j]$ is set to 1, then we know that prefix $P[0..j]$ matches $T[i - j + 1..i]$, and Operation 4 shifts the bits in $B$ by one position, so that in the next iteration we will check whether $P[j + 1]$ matches $T[i + 1]$.

In labeled DAG $G = (V,E)$, each node $v_i \in V$ has a single-character label $\ell(v_i)$. We generalize the shift-and algorithm to labeled DAGs by computing a bit-vector $B_i$ for each node $v \in V$, initializing them to zero. Consider a BFS visit of DAG $G$. When visiting node $v_i$, each bit-vector $B_k$ of its in-neighbour $v_k \in \text{in}(v_i)$ represents a set of prefixes of $P$ matching a path in the graph ending at $v_k$. Thus, we merge all of this information together by taking the bit-wise or of all of the in-neighbours of $v_i$, that is we replace Operation 1 with $B_i \leftarrow 1 + \bigvee_{v_k \in \text{in}(v_i)} B_k$. Operations 2, 3 and 4 are performed as before. An example of the state of the data structures after the execution of the algorithm is shown in Figure 2, and the body of the iteration now is:

1. $B_i \leftarrow 1 + \bigvee_{v_k \in \text{in}(v_i)} B_k$;
2. $B_i \leftarrow B_i \land M[*][T[i]]$;
3. if $B_i[m - 1] = 1$, return yes;
4. $B_i \leftarrow B_i << 1$. 


The adaptation of the classical algorithm for matching pattern $P$ in level DAG $G$. Each bit-vector $D_v$ represent the result after the merging of the bit-vectors of the in-neighbours of $v$ and before the shifting.

5.3 Quantum Bit-Parallel Algorithm for Level DAGs

We make the classic techniques work in a quantum setting for a special class of DAGs, which we call level DAGs. A level DAG is a DAG such that, for every two nodes $v$ and $w$, every path from $v$ to $w$ has the same length, as for the DAG in Figure 2. We also note that degenerate strings [1] can be represented as level DAGs. We use a function representing in-neighbours:

$$in_i(d) = \text{index of the } d\text{-th in-neighbour of } v_i$$

Our approach aims to represent each bit vector $B_i$ with a single qubit $V_i$ set up in a proper superposition, and translate the bit-wise operations to parallel operations across such superposition. In the algorithm, we use the following qubits and quantum registers. Quantum registers $I$ and $J$ store the index of a node and the position in the pattern, respectively. Qubit $V_i$ represents, in superposition, the bit-vector of the node $v_i$, and qubit $E_{i,d}$ stores the contribution of edge $(v_{in_i(d)}, v_i) \in E$ in the update of qubit $V_i$, for, $0 \leq i \leq n - 1$, $0 \leq d \leq D_i - 1$ and $D_i = \text{indeg}(v_i)$. Quantum register $C$ stores label $\ell(v_i)$ of the node in the current iteration, and is used to fetch the content of the corresponding matrix column, which we will store in qubit $M$. Occurrences of the pattern encountered during the execution of the algorithm are stored in qubit $R_i$. Qubits $V_i'$ and $R_i'$ are auxiliary qubits used to store intermediate results, and we also use auxiliary qubits $A$ and $B$ and auxiliary quantum register $Q$ to implement necessary operations. Moreover, we assume to have access to QRAM.

5.3.1 The algorithm

Assume all the quantum registers and qubits to be initialized to $|0\rangle$, except $Q$ initialized to $|1\rangle$. The algorithm starts by setting quantum register $J$ in a balanced superposition, by applying the Hadamard gate on each one of its qubits. Then, we initialize qubits $A$ so that $|a_j\rangle_A = |1\rangle$ for $j = 0$, and $|a_j\rangle_A = |0\rangle$ otherwise. We do the same with qubit $B$, with the difference that $|b_j\rangle_A = |1\rangle$ for $j = m - 1$, and $|b_j\rangle_A = |0\rangle$ otherwise. We can do these operations with two applications of a generalized Toffoli gate, using register $J$ as control and qubits $A$ and then $B$ as targets. In the case of qubit $A$, we first apply an $X$ gate to every qubit of register $J$, we then apply the Toffoli gate, and finally we undo the applications of
We compute the value of \( \ell(\text{node}) \) with the same operations described below for the main loop, the only difference being that the bit-wise or be such that

\[
\text{neighbours, where } \text{node}.
\]

Once we processed the last in-neighbour, \( \text{node} \) is the number of in-neighbours of node \( v_i \).

The rest of the algorithm maintains almost the same overall structure, with the exception of one necessary adaptation. In a DAG of \( L \) levels where \( L_i \) is the set of nodes at level \( l \), for \( 0 \leq l \leq L - 1 \), we iterate over them one at the time, and for each level we process its nodes one after the other. As we will better explain later, we wait before applying the quantum equivalent of the shift operation once we scanned the whole level, not after processing every node. The overall idea is to translate the classical bit-parallel operations into analogous quantum operations that work across the superposition. This translation of bit-parallelism to superposition parallelism is the core of our technique, and we now describe how to apply it to each operation. The pseudocode of the entire procedure is given in Algorithm 1, where all the arithmetic operations are to be considered modulo \( 2^{|P|} \). We only omit the pseudocode for procedures \( \text{SourceNodesInit}() \), \( \text{IncreaseI}() \) and \( \text{IncreaseJ}() \), which is to be found in Appendix B. We also assume \( |P| \) to be a power of two. If this is not the case, we generate a superposition as large as the first power of two greater than \(|P|\), then standard techniques can be used to handle the additional substates, as explained in Appendix A.

**Operation 1 (line 10)** can be broken down into two simpler operations: computing the bit-wise or and adding 1. In our translation to quantum computing, each sub-state of superposition \( \sum_{j=0}^{m-1} |j\rangle_j |v_{i,j}\rangle \) represents an entry of the classical bit-vector used in the Shift-And algorithm. Thus, what was a bit-wise or is now easily translated into the application of few quantum gates. Notice that, to compute the logical or between two generic qubits \( P \) and \( Q \) and store the result in qubit \( R \), we can follow De Morgan’s rules and apply an \( X \) gate to both \( P \) and \( Q \), apply a Toffoli gate with controls \( P \) and \( Q \) and target \( R \), apply an \( X \) gate to \( R \), and finally apply an \( X \) gate to \( P \) and \( Q \) again to restore their initial values. In our case, at iteration \( i \), we use qubit \( E_{i,d} \) to store the \( or \) computed among the first \( d + 1 \) in-neighbours \( v_{i,m(0)}, \ldots, v_{i,m(d)} \) of node \( v_i \), and we compute it in the following way. Let

\[
\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_j |v_{i,m(0)},j\rangle |v_{i,m(1),j}\rangle |v_{i,m(2),j}\rangle \cdots |v_{i,m(d-1),j}\rangle |v_{i,m(d),j}\rangle |\epsilon_{i,d,j}\rangle E_{i,d-1}
\]

be such that

\[
\epsilon_{i,d,j} = v_{i,m(0),j} \lor v_{i,m(1),j} \lor \cdots \lor v_{i,m(d-1),j}.
\]

We compute the value of \( E_{i,d} \) from \( E_{i,d-1} \) and \( V_{i,m(d)} \) as

\[
\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_j |v_{i,m(d),j}\rangle |v_{i,m(d)}\rangle |\epsilon_{i,d-1,j}\rangle E_{i,d-1} \rangle |0\rangle_{E_{i,d}} \rightarrow
\]

\[
\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_j |v_{i,m(d),j}\rangle |v_{i,m(d)}\rangle |\epsilon_{i,d-1,j}\rangle E_{i,d-1} \rangle |v_{i,m(d),j} \lor \epsilon_{i,d-1,j}\rangle E_{i,d}.
\]

Once we processed the last in-neighbour, \( E_{i,D_i-1} \) stores the \( or \) computed among all in-neighbours, where \( D_i \) is the number of in-neighbours of node \( v_i \).
Then, using the same technique as in Operation 1, we compute a \( k \) value. Sub-states are those such that \( |\alpha_j\rangle_A = |\delta_{0,j}\rangle \), we obtain \( |0\rangle_{V_i'} \rightarrow |v'_{i,j}\rangle_{V_i'} \) where \( |v'_{i,j}\rangle_{V_i'} = |1\rangle \) for \( j = 0 \), while \( |v'_{i,j}\rangle_{V_i'} = |e_{i,D_{i-1,j}}\rangle \) for \( 1 \leq j \leq m - 1 \).

**Operation 2 (line 11)** is implemented as a Toffoli-gate application with qubits \( M \) and \( V_i' \) as control and \( V_{i} \) as target.

\[
\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |m_j,\ell(v_i)\rangle_M |v'_{i,j}\rangle_{V_i'} |0\rangle_{V_i} \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |m_j,\ell(v_i)\rangle_M |v'_{i,j}\rangle_{V_i'} |m_j,\ell(v_i) \land v'_{i,j}\rangle_{V_i}
\]

**Operation 3 (line 12)** is replaced by storing in register \( R_i \) the presence of a match ending at node \( v_i \). This requires an intermediate step in which we use qubit \( B \) to filter the content of \( V_{i} \). In fact, qubit \( V_{i} \) now is in state \( |v_{i,j}\rangle_{V_i} = |1\rangle \) for those values of \( j \) such that \( P[0,j] \) has a match ending at \( v_i \) in the graph, and \( |v_{i,j}\rangle_{V_i} = |0\rangle \) otherwise. Since we only care about potential full matches represented by \( |v_{i,m-1}\rangle_{V_i} \), we use \( B \), which is in state \( |\delta_{m-1,j}\rangle_B \), as control qubit of a Toffoli gate, the other control qubit being \( V_{i} \) and the target qubit being \( R'_{i} \).

\[
\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{i,j}\rangle_{V_i} |\delta_{m-1,j}\rangle_B |0\rangle_{R_i'} \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{i,j}\rangle_{V_i} |\delta_{m-1,j}\rangle_B |v_{i,j} \land \delta_{m-1,j}\rangle_{R_{i}'}
\]

Then, using the same technique as in Operation 1, we compute an or between \( R'_i \) and \( R_{i-1} \), storing the result in \( R_i \).

\[
\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{i,j} \land \delta_{m-1,j}\rangle_{R'_i} |r_{i-1,j}\rangle_{R_{i-1}} |0\rangle_{R_i} \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{i,j} \land \delta_{m-1,j}\rangle_{R'_i} |r_{i-1,j}\rangle_{R_{i-1}} |(v_{i,j} \land \delta_{m-1,j}) \lor r_{i-1,j}\rangle_{R_i}
\]

After this operation, \( |r_{i,m-1}\rangle_{R_i} \) is turned to \( |1\rangle \) if there is a full match of \( P \) ending at \( v_i \), otherwise \( |r_{i,m-1}\rangle_{R_i} \) is left unaltered.

**Operation 4 (line 14)** consists in shifting all bits of the classical bit-vector by one position. In the quantum setting, we can perform this operation by adding 1 to index register \( J \) and then reorganising the sum: \( |1\rangle_{C_J} |j\rangle_{J} \rightarrow |1\rangle_{C_J} |j + 1\rangle_{J} \). Notice that this changes value \( |j\rangle_j \), in every term of the superposition to \( |j + 1\rangle_j \). This can be interpreted as “shifting” value \( k_j \) of generic register \( K \) from \( |j\rangle_J |k_j\rangle_K \) to \( |j + 1\rangle_J |k_j\rangle_K \). Because this operation acts on every quantum register and qubit in this way, we have to reset qubits \( A \) and \( B \) to \( |0\rangle \) before performing this operation and reinitialize their values afterwards, so that we prevent their values to be shifted. For the same reason, we also have to wait until having processed the whole level, otherwise we would shift the values of all the nodes at the previous level and compromise the computation.

As last step of the algorithm, we run Grover’s search that uses as oracle function the whole procedure described up to this point, and then applies a Z gate on qubit \( R \). Thus, the marked sub-states are those such that \( |r_{i,j}\rangle_{R_i} = |1\rangle \), which get mapped to \( |r_{i,j}\rangle_{R_i} \). Sub-states such that \( |r_{i,j}\rangle_{R_i} = |0\rangle \) remain unaltered. As for the case of string matching in plain text, we rerun the whole algorithm a constant number of times to boost the probability of success, as explained in Theorem 5 and Appendix C. Algorithm 1 shows the entire procedure.

To prove the correctness of Algorithm 1, we formalise the key properties in the following lemmas. We start by ensuring that the shift operation provides the desired result. Let \( l \) and \( y \) be the total number of times that we started the execution of the middle for-loop
Algorithm 2 Algorithm for testing whether pattern string $P$ has a match in level DAG $G$, running in $O(|E|/\sqrt{|P|})$.

1. time.

Input: Graph $G$, pattern $P$, and constant $c$.

Output: Returns yes if $P$ occurs in $G$, otherwise no.

2. for $c$ times do
3. Initialize quantum register $Q$ to $|1\rangle$;
4. Initialize quantum registers $I,J,A,B,C,M,V_0,V_1,...,V_{n-1},V'_0,...,V'_{n-1}$,
   where $i \in [0,n-1]$ and $d \in [0,D_i-1]$;
   // Apply Hadamard to $J$
5. $|0\rangle_J \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J$;
6. SourceNodesInit($I,J,A,C,V_0,\ldots,V_{n-1},V'_0,\ldots,V'_{n-1}$);
7. IncreaseJ($J,A,B$);
   // $L$ is the number of levels
8. for $l \in [1,L-1]$ do // scan every level
   9. for $|L_l|$ times do // scan every node in the level
10. OperationOne($l,I,C,M,E_i,0,\ldots,E_i,D_i-1,V'_i$);
11. OperationTwo($M,V'_i,V_i$);
   // Invariant 1 holds here
12. OperationThree($B,V_i,R'_i,R_i$);
13. IncreaseI($I,M,C$);
14. OperationFour($J,A,B$);
   // Invariant 2 holds here
15. Apply gate $Z$ to qubit $R_{n-1}$, so that the sign of the amplitude is flipped if
   $|r_{n-1,j}\rangle_{R_{n-1}} = |1\rangle$;
16. Choose $K \in [0,|P|]$ uniformly at random;
17. Run Grover’s iterate operator the optimal number of times assuming to have $K$
   solutions, with the oracle function being lines 6–15 of this algorithm;
18. Measure $R_{n-1}$ into classical register $R_d$;
19. if $R_d = 1$ then
20. return yes
21. return no
Function OperationOne($l, I, C, M, E_{i,0}, ..., E_{i,D_i-1}, V'_i$):
1 for $|L_i|$ times do // scan every node in the level
2 // scan every node in $in(v_i)$
3 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |i\rangle_I |0\rangle_C \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |i\rangle_I |\ell(v_i)\rangle_C; \]
4 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |\ell(v_i)\rangle_C |0\rangle_M \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |\ell(v_i)\rangle_C |m_{\ell(v_i),j}\rangle_M; \]
5 \[ k \leftarrow in_i(0); \quad \text{// Classical operation} \]
6 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{k,j}\rangle_{V_k} |0\rangle_{E_i,0} \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{k,j}\rangle_{V_k} |v_{k,j}\rangle_{E_i,a}; \]
7 for $d \in [1, D_i - 1]$ do // scan every node in $in(v_i)$
8 \[ k \leftarrow in_i(d); \quad \text{// Classical operation} \]
9 // Add the contribution of the current in-neighbour
10 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{k,j}\rangle_{V_k} |\ell_{i,d-1,j}\rangle_{E_i,d-1} |0\rangle_{E_i,d} \rightarrow \]
11 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{k,j}\rangle_{V_k} |\ell_{i,d-1,j} \lor v_{k,j}\rangle_{E_i,a}; \]
12 // Turn to $|1\rangle$ the substate corresponding to $j = 0$
13 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |\delta_{0,j}\rangle_A |\ell_{i,D_i-1,j} \lor v_{k,j}\rangle_{E_i,D_i-1} |0\rangle_{V'_i} \rightarrow \]
14 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |\delta_{0,j}\rangle_A |\ell_{i,D_i-1,j} \lor v_{k,j}\rangle_{E_i,D_i-1} |v_{k,j} \lor \delta_{0,j}\rangle_{V'_i}; \]

Function OperationTwo($M, V'_i, V_i$):
15 // Compute the and with the column of the matrix
16 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |m_{\ell(v_i),j}\rangle_M |v'_{i,j}\rangle_{V'_i} |0\rangle_{V_i} \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |m_{\ell(v_i),j}\rangle_M |v'_{i,j}\rangle_{V'_i} |m_{\ell(v_i),j} \land v'_{i,j}\rangle_{V_i}; \]

Function OperationThree($B, V_i, R'_i, R_i$):
17 // Set $|v_{i,m-1}\rangle_{R_i} = |1\rangle$ if there is a match ending at $v_i$
18 // Apply Toffoli on $V_i$, $B$ and $R'_i$
19 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{i,j}\rangle_{V_i} |\delta_{m-1,j}\rangle_B |0\rangle_{R'_i} \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{i,j}\rangle_{V_i} |\delta_{m-1,j}\rangle_B |v_{i,j} \land \delta_{m-1,j}\rangle_{R'_i}; \]
20 // Apply logic or on $R'_i$, $R_i-1$ and $R_i$
21 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{i,j} \land \delta_{m-1,j}\rangle_{R_i} |r_{i-1,j}\rangle_{R_i-1} |0\rangle_{R_i} \rightarrow \]
22 \[ \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |v_{i,j} \land \delta_{m-1,j}\rangle_{R_i} |r_{i-1,j}\rangle_{R_i-1} |(v_{i,j} \land \delta_{m-1,j}) \lor r_{i-1,j}\rangle_{R_i}; \]

Function OperationFour($J, A, B$):
23 \[ \text{IncreaseJ}(J, A, B); \]
(lines 9–13) and of the outer for-loop (lines 8–14), respectively. That is, \( y = x + \sum_{\lambda=1}^{l-1} |L_\lambda| \) for \( l \geq 2 \), where \( x \in [0, |L_0|] \) is the number of times that we started the execution of the middle for-loop during the \( l \)-th iteration of the outer for-loop. Notice that \( y = 0 \) when \( l = 0 \), and \( y = x \) when \( l = 1 \).

**Lemma 3 (Invariant 1).** During the \( l \)-th execution of the outer for-loop (lines 8–14) and the \( y \)-th execution of the middle for-loop (lines 9–13) of Algorithm 1, but before the \( y \)-th execution of OperationThree() (line 14), Invariant 1 holds: for every qubit \( V_i \), such that \( i \in L_1 \) and \( i \leq t \), we have \( |v_{i,j}|_{V_i} = |1\rangle \) if and only if there exists a path in \( G \) ending at \( v_i \) and matching \( P[0,j] \), where \( t = |L_0| + x - 1 \) is the index of the last node \( v_t \) Algorithm 1 visited so far.

**Proof.** We proceed by strong induction on \( y \), defined as above.

**Base case, \( y = 0 \).** In this case, we executed the initialization but we have not run yet neither the outer nor the middle for-loop. Thus, \( l = 0 \), \( t = |L_0| - 1 \), and qubits \( V_i \) such that \( i \in L_0 \) and \( i \leq t \) are those with in-degree zero, which are initialized by function SourceNodesInit(). For each such \( i \), given that \( J \) is in state \( \sum_{j=0}^{m-1} |j\rangle \langle j| \), function SourceNodesInit() first loads character \( \ell(v_i) \) in register \( C \) and matrix entry \( m_{\ell(v_i),j} \) in register \( M \), in superposition. Then, with regard to \( t \), it performs transformation

\[
\sum_{j=0}^{m-1} |m_{\ell(v_i),j} \rangle_M |\delta_{0,j} \rangle_A |0\rangle_{V_i} \rightarrow \sum_{j=0}^{m-1} |m_{\ell(v_i),j} \rangle_M |\delta_{0,j} \rangle_A |m_{\ell(v_i),j} \wedge \delta_{0,j} \rangle_{V_i},
\]

where, by definition, \( v_{i,j} = m_{\ell(v_i),j} \wedge \delta_{0,j} \). Thus, \( |v_{i,j}| = |0\rangle \) for every \( j \neq 0 \) because of \( \delta_{0,j} \), and \( |v_{i,j}|_{V_i} = |m_{\ell(v_i),j} \rangle_{V_i} \) for \( j = 0 \), which in turn means that \( |v_{i,0}|_{V_i} = |1\rangle \) if and only if \( P[0,0] = \ell(v_i) \).

**Inductive case, \( y \geq 1 \).** We further divide our analysis in two sub-cases.

**First sub-case, \( x = |L_1| \).** In this case, \( y \) is the last iteration of the inner for-loop during the \( l \)-th iteration of the outer for-loop. We assume the inductive hypothesis to hold after the execution of OperationTwo(). We execute OperationThree() and IncreaseI(), which do not change the state of any \( V_z \), for any \( z \in [0, |V| - 1] \). Now, we have to perform OperationFour() (line 14) before starting iteration \( y + 1 \) of the middle for-loop, which will start iteration \( l + 1 \) of the outer for-loop. Assuming the inductive hypothesis, the application of OperationFour() makes every \( V_z \) with \( z \in L_1 \) such that \( |j'\rangle_j |v_{z,j}\rangle_{V_z} = |j'\rangle |1\rangle \), where \( j' = j + 1 \), if and only if there is a match for \( P[0,j] \) in \( G \) ending at \( v_z \), otherwise \( |j'\rangle_j |v_{z,j}\rangle_{V_z} = |j'\rangle |0\rangle \). Then, we start iteration \( y + 1 \) (\( l + 1 \)). Notice that we update \( V_i \) if and only if \( i \in L_{t+1} \) and, in any previous iteration of the middle for-loop, this could have never been the case, thus every \( |v_{i,j}|_{V_i}, i \in L_{t+1} \), is currently set to \( |0\rangle \). The same holds for every \( V'_z \). The for-loop inside OperationOne() computes a logic \( or \) between all the qubits representing all the in-neighbours of \( v_i \). Indeed, before running this for-loop, we have \( |j'\rangle_j |v_{i,m(0),j}\rangle_{E_i^0} \). After one iteration, we have \( |j'\rangle_j |v_{i,m(0),j} \lor v_{i,m(1),j}\rangle_{E_i^1} \). After two iteration, we have \( |j'\rangle_j |v_{i,m(0),j} \lor v_{i,m(1),j} \lor v_{i,m(2),j}\rangle_{E_i^2} \). After \( D_1 - 1 \) iterations, we have \( |j'\rangle |e_{i,D_1-1,j'}\rangle_{E_i,D_1-1} \), where

\[
e_{i,D_1-1,j'} = \bigvee_{d=0}^{D_1-1} v_{i,m(d),j}.
\]
We store an intermediate result in \( V'_i \), \( |v'_{i,j'}\rangle_{V'_i} \), where \( v'_{i,j'} = e_i, D_{i-1,j'} \) except for \( j' = 0 \), because we make sure that \( |v'_{i,0}\rangle_{V'_i} = |1\rangle \) thanks to the or operation with qubit \( A \), which stores \( |\delta_{0,j'}\rangle_A \). Now we compute the logical and with the entry of the matrix, as in the base case, obtaining \( |v_{i,j'}\rangle_{V_i} \), where

\[
v_{i,j'} = m_{\ell(v_i),j'} \land (\delta_{0,j'} \lor \bigwedge_{d=0}^{D_{i-1}} v_{m_i(d),j}).
\]

Applying the inductive hypothesis, this translates to

\[
v_{i,j+1} = (P[j+1] = \ell(v_i)) \land \bigwedge_{d=0}^{D_{i-1}} P[0,j] \text{ has a match ending at } v_{m_i(d)}
\]

Thus, the statement of the lemma holds for \( y + 1 \).

**Second sub-case, \( x < |L_i| \).** The reasoning is analogous to the previous case, the only difference being that \( j \) does not increase and thus we have to look back by \( x + 1 \) iterations, when \( j \) was increased the last time. This requires to assume that the inductive hypothesis was holding for iteration \( y - x \), that is correct because, by strong induction, we assume the inductive hypothesis to hold for every \( y' \leq y \) while proving the statement for \( y + 1 \).

▶ **Lemma 4 (Invariant 2).** After line 14 of Algorithm 1, Invariant 2 holds: if there exists at least one match for \( P \) in \( G \) ending at some \( v_i \) such that \( i \leq t \), then there exists at least one \( j, 0 \leq j \leq m - 1 \), such that \( |r_{i,j}\rangle_{R_i} = |1\rangle \), where \( v_t \) is the last node we visited in Algorithm 1 before line 14.

**Proof.** We proceed by induction on the number \( l \) of times that we run the for-loop at lines 8–14.

**Base case, \( l = 0 \).** In this case, nodes \( v_l \) such that \( t \in L_l, l' \leq 0 \) are those with in-degree zero, while the for-loop at lines 8–14 has never run. Since we are visiting only single-node paths and we are assuming that pattern \( P \) has length at least two, there can be no match for \( P \) ending at these nodes. Correctly, \( |r_{i,j}\rangle_{R_i} = |0\rangle \) for every \( 0 \leq j \leq m - 1 \).

**Inductive case, \( 1 \leq l \leq L - 1 \).** By inductive hypothesis, we assume the statement of the lemma to be true right after running iteration \( l \) of the for-loop at lines 8–14, and thus right before executing \( \text{IncreaseJ}() \) at line 14. After the execution of \( \text{IncreaseJ}() \), the new state is \( \sum_{j' = 0}^{m-1} |j'\rangle_{R_i} |r_{i,j'}\rangle_{R_i} \), where \( j' = j + 1 \) and \( v_i \) is the last node visited so far. Then, we start iteration \( l + 1 \), processing \( i' \in L_{l+1}, i' = t + 1 \). We execute \( \text{OperationOne()} \) \( \text{OperationTwo()} \), which do not affect register \( R_{i'} \). Then we run the operations at lines 12–12, obtaining \( |r_{i',j'}\rangle_{R_i} \) where \( r_{i',j'} = (v_{i',j} \land \delta_{m-1,j'}) \lor r_{t,j} \). Let us consider the first time we run the middle for-loop during iteration \( l+1 \) of the outer for-loop. If \( P \) has a match ending at some \( v_{i}, z < i' \), the inductive hypothesis guarantees \( r_{t,j} = 1 \) for some \( j \). Otherwise, if \( P \) does not have any such match, then \( r_{t,j} = 0 \) for all \( j \). In this second case, if \( P \) has a match ending at \( v_{i'} \), we know by Lemma 3 that \( v_{i',m-1} = 1 \). This, combined with the fact that \( \delta_{m-1,m-1} = 1 \), correctly implies that \( r_{i',m-1} = 1 \), proving the statement for this specific \( i' \) and \( j' = m - 1 \). If \( P \) has no match ending at \( v_{i'} \), then \( v_{i',m-1} = 0 \) and \( r_{i',j'} = 0 \) for all \( j' \), which must be the case when no match has been found yet. To conclude the proof, notice that the same reasoning applies for the subsequent iterations of the middle for-loop by using every time the
previous instance of this reasoning in place of the inductive hypothesis. That is, we use \( r_t, j' \)
when proving the statement for \( r_{t+1}, j' \) and so on, until we prove the statement for \( r_1, j' \),
where \( r_t \) is the last node with index in \( L_t+1 \). At this point, we exit the middle \( \text{for} \)-loop and
the statement of the lemma is proven for \( l + 1 \).

The correctness of the algorithm follows from the previous lemma combined with few
additional observations.

\[ \textbf{Theorem 5.} \quad \text{Given pattern string } P \text{ of length at least 2 and level DAG } G, \text{ Algorithm 1}
\] returns the right answer for the SMLG problem on \( P \) and \( G \) with probability \( p > 1 - (7/8)^c \),
for any given integer \( c \).

\[ \textbf{Proof.} \quad \text{After running the outer } \text{for} \)-loop of Algorithm 1 \( L - 1 \) times, we exit such a loop, and
\] we know we have visited all the nodes (nodes in \( L_0 \) where visited during the initialization).
If we consider Lemma 4 applied in the case of \( t = n - 1 \), we are considering all the nodes,
which means that if \( P \) has no match ending in \( G \), then no substate of register \( R_{n-1} \) is such
that \( |r_{n-1, j}\rangle_{R_{n-1}} = |1\rangle \), for any \( j \). Instead, if \( P \) has a match in \( G \), then at least one substate
of \( R_{n-1} \) is such that \( |r_{n-1, j}\rangle_{R_{n-1}} = |1\rangle \), for some \( j \). We use standard techniques that consist
in rerunning the algorithm a constant number of times to boost the probability of measuring
such a state, and achieve the desired one. Appendix C provides a more detailed analysis.

Finally, the time complexity of our algorithm is subquadratic in the size of the graph.

\[ \textbf{Theorem 6.} \quad \text{The time complexity of Algorithm 1 is } O(|E|^{\sqrt{|P|}}) \text{ in the QRAM model, and}
\] the space complexity is \( O(|E| + |V|) \).

\[ \textbf{Proof.} \quad \text{The algorithm uses } |V| \text{ qubits } V_i, \text{ and the same amount of qubits } V'_i, R_i, R'_i; \text{ qubits}
\] \( E_{i, d} \) are a total of \( |E| \) qubits, and the rest are a constant number of qubits and registers.
Thus, the space complexity is \( O(|E| + |V|) \).

With the \( \text{for} \)-loop in function \texttt{SourceNodesInit}(), the algorithm visits the nodes in \( L_0 \),
which are at most \( O(|V|) \). The iteration conditions at lines 8 and 9 make the algorithm visit
every node. For each such iteration, we perform a constant number of operations except
for the \( \text{for} \)-loop in \texttt{OperationOne}(). This \( \text{for} \)-loop visits all the in-neighbours of a node,
each time performing a constant number of operations, and \( \sum_{i=0}^{V_i-1} |in(i)| = |E| \). All of the
aforementioned operations can be implemented with a constant number of quantum-gate
applications, each affecting a constant number of qubits (3 at most), or by performing a load
operation from the QRAM, assumed to require constant time. At the end of the algorithm,
we run Grover’s search procedure on a superposition of \( 2|P| \) states, using the entire algorithm
as the oracle function.

Summing everything up, we spend \( O(|V|) \) time for the initialization, \( O(|V| + |E|) \) time in
the \( \text{for} \)-loops, and \( O(|E|^{\sqrt{|P|}}) \) time for Grover’s search procedure. The total time complexity
is thus dominated by \( O(|E|^{\sqrt{|P|}}) \).

\[ \text{References}
\]
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A Reductions to the power-of-two case

For string matching in plain text, if $|T|$ is not a power of two, in addition to quantum register $I$ for indexing, we use also quantum register $I'$, of the same size. Let $x$ be the only integer $x$ such that $|T| < 2^x < 2|T|$. Generate superposition $\sum_{i=0}^{2^x-1} |i\rangle_I |0\rangle_I$ and compute $\sum_{i=0}^{2^x-1} |i\rangle_I |i' \mod |T|\rangle_I$. Now run the algorithm using $I$ as normal. This creates some redundant substates, but does not affect the correctness of the algorithm.

For SMLG in level DAGs, if $|P|$ is not a power of two, we generate a superposition of size $2^x$, where $x$ is the only integer such that $|P| < 2^x < 2|P|$. Then, it suffices to assume that every entry that we read from the QRAM to the additional substates between $|P|$ and $2^x$.
is always initialized to $|1\rangle$, because this is the neutral value in a logical and, an thus in the application of the Toffoli gate. Therefore, in these substates, a qubit $R_i$ can and will be set to value $|1\rangle_{R_i}$ if and only if a previous “shift” carried $|1\rangle_{R_{i-1}}$.

Alternatively, if $|P|$ is not a power of two, we can classically reduce the problem to this case. We add new symbol $\$ to the alphabet. Then, we pad $P$ with as many $\$ at the end as needed to reach the next power of two. For each level in the DAG, we add a new node with label $\$, and we place an edge for every node in that level to the new node. We connect all this new nodes in a chain, and we also add a chain of $|P|$ such nodes after the last level (they create new levels consisting only of one node). The pattern now can overflow in these nodes after finding a proper match in the DAG. Finally, we apply the same binary encoding as in the plain text case, now replacing every node with a chain of two nodes, sending all the incoming edges to the first node and making all the outgoing edges leave from the second node. Overall, we add one new node per level, and one new edge per node, plus $|P|$ additional nodes and edges after the last level. This takes time $O(|E| + |P|)$.

## B Additional pseudo-code

```plaintext
1 Function SourceNodesInit(I, J, A, C, V_0, ..., V_{n-1}, \ldots, V_{n-1}):
   // Initialize $|a_i\rangle_A$ so that $a_j = 1$ if $j = 0$, $a_j = 0$ otherwise
2 $\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |0\rangle_A \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |\delta_{0,j}\rangle_A$;
   // Initialize $|b_j\rangle_B$ so that $b_j = 1$ if $j = m - 1$, $b_j = 0$ otherwise
3 $\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |0\rangle_B \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |\delta_{m-1,j}\rangle_B$;
   // $L_0$ is the set of nodes in level 0.
4 for $|L_0|$ times do
   // Read node label $\ell(v_i)$ in $C$
5 $\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |i\rangle_I |0\rangle_C \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |i\rangle_I |\ell(v_i)\rangle_C$;
   // Read the matrix entries for character $\ell(v_i)$ in $M$
6 $\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |\ell(v_i)\rangle_C |0\rangle_M \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |\ell(v_i)\rangle_C |m_{\ell(v_i),j}\rangle_M$;
   // Apply Toffoli to qubits $M$, $A$ and $V_i$
7 $\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |m_{\ell(v_i),j}\rangle_M |\delta_{0,j}\rangle_A |0\rangle_{V_i} \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |m_{\ell(v_i),j}\rangle_M |\delta_{0,j}\rangle_A |m_{\ell(v_i),j} \oplus \delta_{0,j}\rangle_{V_i}$;
   // Reset $M$ and $C$
8 $\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |\ell(v_i)\rangle_C |m_{\ell(v_i),j}\rangle_M \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |\ell(v_i)\rangle_C |m_{\ell(v_i),j} \oplus m_{\ell(v_i),j}\rangle_M$;
9 $\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |i\rangle_I |\ell(v_i)\rangle_C \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |j\rangle_J |\ell(v_i)\rangle_C = |1\rangle_Q |i\rangle_I$;
   // Increase $I$ by one to visit the next node
10 $\frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |1\rangle_Q |i\rangle_I \rightarrow \frac{1}{\sqrt{m}} \sum_{j=0}^{m-1} |1\rangle_Q |i + 1\rangle_I$;
```
which means that if we know we have visited all the nodes (nodes in $c$ times, the overall probability of failure (measuring a wrong result) measuring a wrong result is greater than $K$ iterations has period of good solutions and $R$ that $|r_{n-1,j}\rangle_{R_{n-1}} = |1\rangle$, for any $j$. Instead, if $P$ has a match in $G$, then at least one substate of $R_{n-1}$ is such that $|r_{n-1,j}\rangle_{R_{n-1}} = |1\rangle$, for some $j$.

The for loop that we run at the end of the algorithm ensures to achieve high probability of success. The probability of success $p$ in Grover’s search algorithm is the sinusoidal function $p(K) = \sin^2((2K + 1)\theta)$ [5], where $\theta = \sin^{-1}\left(\frac{\sqrt{N}}{M}\right)$. $N$ is the search space, $M$ is the number of good solutions and $K$ is the number of iterations of the Grover’s operator. This function has period $\lambda_M \approx \frac{\pi}{2} \sqrt{\frac{N}{M}} - 1$. Consider the case $M = 1$. If we choose a random number of iterations $K$ between 1 and $\lambda_1$, we have $p(K) \geq \frac{1}{2}$ with probability $1/2$. This is because half of the material of the function is above the horizontal line of 1/2. When $p(K) \geq 1/2$, the probability of measuring a wrong result is $p_{\text{top}} \leq 1/2$. When $p(K) \geq 1/2$, the probability of measuring a wrong result is greater than 1/2, but anyway $p_{\text{bottom}} \leq 1$. If we run the process $c$ times, the overall probability of failure (measuring a wrong result) $p_f$ is then...
\[ p_f = \left( p_{\text{top}} \frac{1}{2} + p_{\text{bottom}} \frac{1}{2} \right)^c \leq \left( \frac{1}{2} \frac{1}{2} + 1 \cdot \frac{1}{2} \right)^c = \left( \frac{3}{4} \right)^c \]

Thus, the probability of success (measuring a correct result) is \( p_s = 1 - \left( \frac{3}{4} \right)^c \).

In the general case \( 1 < M \leq N \), the period \( \lambda_M \) of function \( p(K) \) is smaller than period \( \lambda_1 \) of the case \( M = 1 \). We can still use the same random number of iterations \( K \) between \( 1 \) and \( \lambda_1 \), as nearly half of the material of the function \( p(K) \) is above the horizontal line of \( 1/2 \): the worst case is when \( \lambda_M \) is little over half of \( \lambda_1 \). In this case we know that \( p(K) \) will be sampled uniformly over half of the range of period \( \lambda_1 \), but the other half may have biased sampling. Namely, the other half of the function might have more material below \( 1/2 \) than above. To have a safe estimate, we assume that the probability of returning the wrong result in the biased case is \( p_{\text{biased}} = 1 \). That is, if we run the process \( c \) times, the overall probability of failure (measuring a wrong result) \( p_f \) is then

\[ p_f = \left( p_{\text{biased}} \frac{1}{2} + \left( p_{\text{top}} \frac{1}{2} + p_{\text{bottom}} \frac{1}{2} \right) \right)^c \leq \left( 1 \cdot \frac{1}{2} + \left( \frac{1}{2} \frac{1}{2} + 1 \cdot \frac{1}{2} \frac{1}{2} \right) \right)^c = \left( \frac{7}{8} \right)^c \]

Thus, the probability of success (measuring a correct result) is \( p_s = 1 - \left( \frac{7}{8} \right)^c \). \( \square \)
On the Impact of Morphisms on BWT-Runs

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Abstract

Morphisms are widely studied combinatorial objects that can be used for generating infinite families of words. In the context of Information theory, injective morphisms are called (variable length) codes. In Data compression, the morphisms, combined with parsing techniques, have been recently used to define new mechanisms to generate repetitive words. Here, we show that the repetitiveness induced by applying a morphism to a word can be captured by a compression scheme based on the Burrows–Wheeler Transform (BWT). In fact, we prove that, differently from other compression-based repetitiveness measures, the measure $r_{\text{bwt}}$ (which counts the number of equal-letter runs produced by applying BWT to a word) strongly depends on the applied morphism. More in detail, we characterize the binary morphisms that preserve the value of $r_{\text{bwt}}(w)$, when applied to any binary word $w$ containing both letters. They are precisely the Sturmian morphisms, which are well-known objects in Combinatorics on words. Moreover, we prove that it is always possible to find a binary morphism that, when applied to any binary word containing both letters, increases the number of BWT-equal letter runs by a given (even) number. In addition, we derive a method for constructing arbitrarily large families of binary words on which BWT produces a given (even) number of new equal-letter runs. Such results are obtained by using a new class of morphisms that we call Thue–Morse-like. Finally, we show that there exist binary morphisms $\mu$ for which it is possible to find words $w$ such that the difference $r_{\text{bwt}}(\mu(w)) - r_{\text{bwt}}(w)$ is arbitrarily large.

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1 Introduction

The Burrows–Wheeler transform (BWT) is a reversible permutation of words, introduced in the Data compression field [5]. Such a transformation allows one to boost the effect of the run-length encoding with respect to the original word in input [27]. Due to its...
myriad virtues, some of the well-known compressed text-indexes for pattern matching [11, 13] and the most used alignment tools in Bioinformatics [23, 21] are based on the BWT. The performance of the BWT is related to the repetitions of factors in the word, which is why the number of equal-letter runs of the BWT, denoted by $r_{bwt}$, is considered as a measure of repetitiveness [29]. Much attention has recently paid to the measure $r_{bwt}$ both for its crucial role in designing compressed indexing data structures for highly repetitive texts [13, 16, 30] and for its combinatorial properties [25, 14].

In Combinatorics on words, morphisms are a fundamental tool for generating repetitive sequences, with multiple applications. For instance, injective morphisms, known as codes, are widely used in the fields of Information theory, Data compression, and Cryptography [2]. Recently, morphisms have been used in conjunction with copy-paste mechanisms to define novel compressors and repetitiveness measures, called NU-systems [32]. Informally speaking, a morphism is a mechanism that transforms each letter in a given input word into a corresponding image word, thus producing an output that is likely to contain longer repeated factors. The relationship between morphisms and the measure $r_{bwt}$ has been studied in the context of a subclass of infinite words generated by morphisms, i.e., the purely morphic words [4, 12].

Here, we focus the impact of morphism application on the number of BWT equal-letter runs of finite words.

In Section 3, we prove that a binary morphism is cyclic (i.e., the images of both letters are powers of the same word) if and only if the image of every word under this morphism has the same number of BWT equal-letter runs, regardless of the input word. We also prove other results relating morphisms and words sharing the same Parikh vector (i.e., having the same number of occurrences of each letter), which can be of independent interest.

Then, in Section 4 we find a novel characterization of Sturmian morphisms [3, 28] in terms of BWT equal-letter runs: they are exactly the binary morphisms that preserve the number of BWT equal-letter runs of every binary word containing both letters of the alphabet. This characterization is interesting from a combinatorial point of view, because Sturmian morphisms are a widely studied subject [3, 28]. It also builds another bridge between Combinatorics on words and Data compression.

Further, in Section 5 we show a wide class of morphisms, which we call Thue–Morse-like morphisms, that increase the number of BWT equal-letter runs by 2 on every binary word containing both letters of the alphabet. Moreover, for each even number $2k$, we can find a wide class of binary morphisms, obtained by composing Sturmian and Thue–Morse-like morphisms, that increase the BWT equal-letter runs of every binary words by exactly $2k$. Note that this is exhaustive for the binary alphabet. In fact, unless considering powers of a single letter, every binary word has an even number of BWT equal-letter runs. In addition, we can use the aforementioned morphisms to construct arbitrarily large families of binary words having all the same number of BWT equal-letter runs, for every fixed (even) number, and converging to an infinite aperiodic word.

At the other end of the spectrum, in Section 6 we show that there are binary morphisms (in particular, the so-called period-doubling morphism) that can highly increase the number of BWT equal-letter runs of binary words. We show that the increase in the number of BWT equal-letter runs can be $\Omega(\sqrt{n})$, where $n$ is the length of the original word. In Section 7, we show that this degree of increase cannot occur in other relevant reachable repetitiveness measures, like the size of the Lempel–Ziv parsing [9, 22], or the size $g$ of the smallest deterministic context-free grammar generating the word [18].

We conclude in Section 8 with some final remarks, and some open questions and conjectures.
2 Preliminaries

Basic terminology

Let $\Sigma = \{a_1, a_2, \ldots, a_\sigma\}$ be a finite sorted set of letters $a_1 < a_2 < \cdots < a_\sigma$, which we call an alphabet. A finite word $w = w[1]w[2] \cdots w[|w|]$ is any finite sequence of letters where $w[i] \in \Sigma$, for $i \in [1, |w|]$, and $|w|$ is the length of the word. We denote by $\text{alph}(w)$ the set of the letters of $\Sigma$ appearing in $w$. The empty word, denoted by $\varepsilon$, is the unique word of length 0. The set of all finite words (resp. all finite words of positive length) over the alphabet $\Sigma$ is denoted by $\Sigma^*$. A factor of $w$ is then called a word of all the $w$ when $w$ of all the unique infinite word $\mu$ of a word is then called a morphism is called a primitive morphism. For some $\mu \in \Sigma^*$ such that $\mu(w[i]) = \mu(w[i+1])$, we denote by $\Pi_{i=1}^{k} w_i$ the concatenation of the words $w_1, w_2, \ldots, w_k$ in that order. We denote by $w^k$ the concatenation of the word $w$ with itself $k$ times. A rotation of the word $w = w[1]w[2] \cdots w[n]$ is a word of the form $w[i+1,n][w[1,i]]$, for some $1 \leq i \leq n$, obtained by shifting $i$ letters cyclically. We denote by $R(w)$ the multiset of all the $w$ rotations of $w$. A factor of any word in $\mathcal{R}(w)$ is called a circular factor of $w$. A word is primitive if $w = w^k$ implies $k = 1$, or equivalently, if it cannot be written as $uv$ for some non-empty words $u$ and $v$ such that $uv = vu$. A primitive word of length $n$ has exactly $n$ distinct rotations. If $w$ is a binary word over the alphabet $\{a, b\}$, the complement of $w$, i.e., the word obtained by replacing all the $a$'s of $w$ by $b$'s and all the $b$'s by $a$'s, is denoted by $\overline{w}$. If $w = w[1] \cdots w[n]$, the reverse of $w$ is the word $w^R = w[n] \cdots w[1]$. Given a word $w \in \Sigma^*$ and $a \in \Sigma$, we denote by $|w|_a$ the number of occurrences of $a$ in $w$. The run-length encoding of a word $w$, denoted by $\text{rle}(w)$, is a sequence of pairs $(c_i, l_i)$ with $c_i \in \Sigma$ and $l_i > 0$, such that $w = c_1^l_1 c_2^l_2 \cdots c_n^l_n$ and $c_i \neq c_{i+1}$. The length $|\text{rle}(w)|$ is the number of equal-letter runs in $w$. The Parikh vector of $w$, denoted as $P(w)$, is the $\sigma$-tuple $(|w|_{a_1}, \ldots, |w|_{a_{\sigma}})$. Given two words $u$ and $v$ having the same length, the Hamming distance between $u$ and $v$, denoted as $d_H(u,v)$, is the number of positions at which the corresponding letters in $u$ and $v$ are different. An infinite word $x = x[1]x[2]x[3] \cdots$ is a non-ending sequence of elements of the alphabet $\Sigma$. An infinite word $x$ is ultimately periodic if there exist $w \in \Sigma^*$ and $v \in \Sigma^*$ such that $x = uvv \cdots$; it is called periodic when $u = \varepsilon$; aperiodic if it is not ultimately periodic. If there is no ambiguity, finite words are simply called words.

Morphisms

Let $\Sigma$ and $\Gamma$ be two alphabets. A morphism is a map $\mu$ from $\Sigma^*$ to $\Gamma^*$ such that $\mu(uv) = \mu(u)\mu(v)$ for all words $u, v \in \Sigma^*$. Therefore, a morphism $\mu$ can be defined by specifying its action on the letters of $\Sigma$ and can be denoted as $\mu \equiv (\mu(a_1), \ldots, \mu(a_\sigma))$. When $\Sigma = \Gamma = \{a, b\}$, $\mu$ is called a binary morphism. A morphism $\mu$ is called prolongable on a letter $a \in \Sigma$ if $\mu(a) = au$ for some $u \in \Sigma^+$. If for all $a \in \Sigma$ it holds that $\mu(a) \neq \varepsilon$, then the morphism $\mu$ is called non-erasing. From now on, we will consider non-erasing morphisms, unless stated explicitly otherwise. If there exists $k$ such that $|\mu(a)| = k$ for every $a \in \Sigma$, then the morphism is called $k$-uniform. A 1-uniform morphism is called a coding. Given a morphism $\mu$ prolongable on some letter $a \in \Sigma$, the family of words $\{a, \mu(a), \mu^2(a), \ldots\}$ are prefixes of a unique infinite word $\mu^{\infty}(a) = \lim_{n \to \infty} \mu^n(a)$, that is a fixed point of $\mu$. Such an infinite word is then called purely morphic. An infinite word is morphic if it is obtained by applying a coding to a purely morphic word. A morphism $\mu$ is cyclic if there exists $z \in \Gamma^*$ such that $\mu(a) \in z^\ast$, for each $a \in \Sigma$. Otherwise, it is called acyclic. Note that the fixed point of a cyclic morphism is periodic. In the case of a binary morphism, it is known that $\mu$ is cyclic if and only if $\mu(ab) = \mu(ba)$.
Sturmian words and Sturmian morphisms

Let $\Sigma = \{a, b\}$. A word $w \in \Sigma^*$ is called balanced if the difference of the number of $a$’s (or, equivalently, $b$’s) in every two factors of the same length of $w$ is at most 1. An infinite word $x$ is balanced if every finite factor of $x$ is balanced. A finite word $w$ is circularly balanced if each word in $R(w)$ is balanced.

An infinite word over $\Sigma = \{a, b\}$ is a Sturmian word if it has exactly $n + 1$ distinct factors of length $n$ for every $n \geq 0$. The theory of Sturmian words is very well studied (see [24] for a reference). For example, the following characterization is well known.

$\triangleright$ Theorem 1. An infinite word over $\Sigma = \{a, b\}$ is Sturmian if and only if it is balanced and aperiodic.

A class of Sturmian words, called characteristic Sturmian words, can be constructed by using finite words, called standard Sturmian words, defined recursively as follows. Given an infinite sequence of integers $(d_0, d_1, d_2, \ldots)$, with $d_0 \geq 0, d_i > 0$ for all $i > 0$, called directive sequence, the associated standard Sturmian words are defined by $s_0 = b, s_1 = a$, and $s_{i+1} = s_{d_i-1}s_{i-1}$, for $i \geq 1$. A characteristic Sturmian word is the limit of an infinite sequence of standard Sturmian words, i.e., $s = \lim_{i \to \infty} s_i$. Note that standard Sturmian words are finite words also appearing as extremal case for several algorithms and data structures [19, 7, 26, 37].

A Sturmian morphism is a morphism that maps infinite Sturmian words to infinite Sturmian words. Some combinatorial characterizations of Sturmian morphisms have been proved in [3]. In particular, a binary morphism $\mu$ is Sturmian if and only if it is acyclic and balanced (i.e., it maps balanced words to balanced words). Berstel and Séébold [3] also proved the following characterization:

$\triangleright$ Theorem 2. An acyclic morphism $\mu$ is Sturmian if and only if it is locally Sturmian, that is, there exists a Sturmian word $s$ such that $\mu(s)$ is Sturmian.

Let us denote the following morphisms:

$$E : \begin{cases} a \mapsto b \\ b \mapsto a \end{cases} \quad \phi : \begin{cases} a \mapsto ab \\ b \mapsto a \end{cases} \quad \tilde{\phi} : \begin{cases} a \mapsto ba \\ b \mapsto a \end{cases}$$

The morphism $\phi$ is called the Fibonacci morphism, since its fixed point is the Fibonacci word $abaababaababaab \cdots$. The monoid $\{E, \phi, \tilde{\phi}\}^*$ generated by $E$, $\phi$, and $\tilde{\phi}$, by using the composition operator $\circ$, is known as the Sturm monoid. The following theorem, proved in [28], shows the combinatorial structure of Sturmian morphisms.

$\triangleright$ Theorem 3. A morphism is Sturmian if and only if it belongs to $\{E, \phi, \tilde{\phi}\}^*$.

Burrows–Wheeler transform

The Burrows–Wheeler transform (BWT) of a word $w$, denoted by $\text{bwt}(w)$, is a permutation of $w$ obtained by sorting all its rotations in lexicographical order and then concatenating the last symbol of each rotation. The original word can be recovered if one stores the position where it appears in the list of sorted rotations. If a word is highly repetitive, the number of equal-letter runs of the BWT tends to be small. In fact, Kempa and Kociumaka have shown that $r_{\text{bwt}}$ is never too far from the size of the Lempel-Ziv parsing, a widely used repetitiveness measure [16]. Hence applying run-length encoding to the BWT is very effective. Because of this, the value $r_{\text{bwt}}(w) = |\text{rle}(\text{bwt}(w))|$ that counts the number of $BWT$-runs of $w$, i.e., equal-letter runs of $\text{bwt}(w)$, is used as a measure for capturing the repetitiveness of the word.
The equivalence of 1, 2 and 3 is in [26, 35]. The equivalence with 4 is in [8] (see also Proposition 10 in [34]).

3 Morphisms and sorted rotations of words

We start by introducing some definitions regarding the rotations of morphic images of words.

Definition 5. Let $\mu : \Sigma^* \to \Gamma^*$ be a morphism. Then, we define the multisets

\[ I_\mu(w) = \{\mu(w') \mid w' \in \mathcal{R}(w)\} \]

\[ S_\mu(w) = \{v\mu(w')u \mid u, v \in \Gamma^+, vw = \mu(a) \text{ for some } a \in \Sigma, \text{ and } au \in \mathcal{R}(w)\}. \]

The multiset $I_\mu(w)$ corresponds to the rotations of $\mu(w)$ obtained by applying $\mu$ to the rotations of $w$. The multiset $S_\mu(w)$ corresponds to all the remaining rotations of $\mu(w)$. We refer to the multiset $I_\mu(w)$ as the I-rotations of $\mu(w)$, and to the multiset $S_\mu(w)$ as the S-rotations of $\mu(w)$. These two multisets could have elements that end up being equal, as we show in the following example.

Example 6. Let $\mu \equiv (a, bab)$, which is an acyclic binary morphism. Then, $ab$ is primitive but $\mu(ab) = abab$ is not. Moreover, $I_\mu(w) = \{abab, baba\} = S_\mu(w)$.

We now prove some combinatorial properties of words having the same Parikh vector. By using such properties, we prove that, in the case of the binary alphabet, the lexicographic order among the rotations of a given word is either preserved or reversed, after a morphism is applied. This is a key point to show that the number of BWT-runs cannot decrease after the application of a binary morphism. This is no longer true for larger alphabets.

The following lemma shows that distinct words having the same Parikh vector must have Hamming distance of at least 2.

Lemma 7. Let $w_1, w_2 \in \Sigma^*$ be such that $w_1 \neq w_2$ and $P(w_1) = P(w_2)$. Then, $d_H(w_1, w_2) \geq 2$.

Proof. By definition of $d_H$, we have that $d_H(w_1, w_2) = 0$ if and only if $w_1 = w_2$. So, let us suppose by contradiction that $d_H(w_1, w_2) = 1$. Then, there exist two finite words $u, v \in \Sigma^*$ and two distinct indices $i < j \in [1, \sigma]$ such that $w_1 = ua_i v$ and $w_2 = ua_j v$. It follows that the Parikh vectors of $w_1$ and $w_2$ are respectively

\[ P(w_1) = (|u|_{a_1} + |v|_{a_1}, \ldots, |u|_{a_i} + |v|_{a_i} + 1, \ldots, |u|_{a_j} + |v|_{a_j}, \ldots, |u|_{a_\sigma} + |v|_{a_\sigma}) \]

and

\[ P(w_2) = (|u|_{a_1} + |v|_{a_1}, \ldots, |u|_{a_i} + |v|_{a_i}, \ldots, |u|_{a_j} + |v|_{a_j} + 1, \ldots, |u|_{a_\sigma} + |v|_{a_\sigma}). \]

Thus, we obtain that the $P(w_1) \neq P(w_2)$, a contradiction. ▶
Since all the words in the same conjugacy class share the same Parikh vector, we can derive the following

\textbf{Corollary 8.} Let \( w \in \Sigma^* \) be a word. Then, for every word \( w' \in R(w) \) such that \( w' \neq w \), one has \( d_H(w, w') \geq 2 \).

Here, we introduce and study new properties of some classes of morphisms, which are related to the number of BWT-runs.

\textbf{Definition 9.} A morphism \( \mu \) is abelian order-preserving if for every pair of distinct words \( x \) and \( y \) having the same Parikh vector, it holds that \( x < y \iff \mu(x) < \mu(y) \).

A morphism \( \mu \) is abelian order-reversing if for every pair of distinct words \( x \) and \( y \) having the same Parikh vector, it holds that \( x < y \iff \mu(x) > \mu(y) \).

In general, a morphism can be neither abelian order-preserving nor abelian order-reversing:

\textbf{Example 10.} A cyclic morphism is trivially not abelian order-preserving nor abelian order-reversing. The acyclic morphism \( \mu \equiv (b, a, c) \) is also neither of them. This can be verified on the rotations of the word \( abc \).

However, all acyclic morphisms with a binary domain are either abelian order-preserving or abelian order-reversing, as we show in the following lemma.

\textbf{Lemma 11.} Let \( \mu : \{a, b\}^* \to \Sigma^* \) be an acyclic morphism. Then, \( \mu \) is either abelian order-preserving or abelian order-reversing.

\textbf{Proof.} Let \( \mu \equiv (\alpha, \beta) \) be an acyclic morphism (i.e., \( \alpha \beta \neq \beta \alpha \)). For the proof, we assume that \( |\alpha| \leq |\beta| \), and the other case is treated symmetrically. Factorize \( \mu \) as \( (\alpha, \beta) = (\alpha, \alpha^k v) \), where \( k \geq 0 \) is as big as possible. This factorization is unique, and \( \alpha \) is not a prefix of \( v \), otherwise, \( k \) is not as big as possible. Also, \( v \neq \varepsilon \) and \( v \neq \alpha \) because the morphism \( \mu \) is acyclic. Let \( x = uaz_1 \) and \( y = ubz_2 \) be two distinct binary words with the same Parikh vector. Note that \( a \) has to appear in \( z_1 \), since otherwise \( x \) has fewer \( b \)’s than \( y \). Let \( z_1 = a^tbz_1' \) for some \( t \geq 0 \) and \( z_1' \in \{a, b\}^* \). We can write \( x = uaa'^tbz_1' \). Then, \( \mu(x) = \mu(u)\alpha^k\alpha^tvu(z_1') \) and \( \mu(y) = \mu(u)\alpha^k\mu(z_2) \). We proceed by case analysis.

If \( v \) is not a prefix of \( \alpha \), then the order between \( \mu(x) \) and \( \mu(y) \) depends only on the order between \( \alpha \) and \( v \). The reason is that \( \mu(x) \) and \( \mu(y) \) share a common prefix \( \mu(u)\alpha^k \), followed by \( \alpha \) and \( v \) respectively, which differ at some position from left to right. Hence, if \( \alpha < v \), we obtain \( x < y \iff \mu(x) < \mu(y) \); if \( v < \alpha \), then we obtain \( x < y \iff \mu(x) > \mu(y) \).

If \( v \) is a proper prefix of \( \alpha \) and \( k > 0 \), rewrite \( \mu(y) = \mu(u)\alpha^kvaz_2' \). We can do this because \( y \) has to have at least one letter after \( ub \) and both images \( \alpha \) and \( \beta \) start with \( \alpha \) (in the case of \( \beta \) because \( k > 0 \)). We note that the common prefix \( \mu(u)\alpha^k \) is followed by \( \alpha v \) in \( \mu(x) \) (\( \alpha v \) is a prefix of \( \alpha \alpha \)), and by \( v \alpha \) in the case of \( \mu(y) \). The order between \( \mu(x) \) and \( \mu(y) \) is then completely determined by the order between \( \alpha v \) and \( v \alpha \). This happens because \( \alpha v \) and \( v \alpha \) are words of the same length which must be distinct, as implied by the inequality \( \alpha \beta = \alpha \alpha^k v \neq \beta \alpha = \alpha^k v \alpha \). Hence, if \( \alpha v < v \alpha \), we obtain \( x < y \iff \mu(x) < \mu(y) \); if \( v \alpha < \alpha v \), then we obtain \( x < y \iff \mu(x) > \mu(y) \).

No other case is possible. By construction, \( \alpha \) is not a prefix of \( v \). Also, \( \alpha \neq v \), so if \( v \) is a prefix of \( \alpha \), it has to be a proper prefix. If this is the case, as \( |\alpha| \leq |\alpha^k v| \) and \( |v| < |\alpha| \), \( k \) has to be at least 1.

Using Lemma 11 we can easily derive the following corollary.
Corollary 12. Let \( w \) be a binary word and let \( \mu \) be an acyclic morphism. Then, for all pairs of rotations \( u, v \) of \( w \), either \( u < v \iff \mu(u) < \mu(v) \) (when \( \mu \) is abelian order-preserving), or \( u < v \iff \mu(u) > \mu(v) \) (when \( \mu \) is abelian order-reversing).

We introduce new measures to study how the action of a morphism affects the BWT-runs.

Definition 13. Let \( \mu \) be a morphism and \( w \) a word. We define
\[
\Delta^+_\mu(w) = r_{bwt}(\mu(w)) - r_{bwt}(w)
\]
and
\[
\Delta^-\mu(w) = \frac{r_{bwt}(\mu(w))}{r_{bwt}(w)}.
\]

Acyclic binary morphisms cannot decrease the number of BWT-runs of any word.

Theorem 14. Let \( \mu : \{a,b\}^* \rightarrow \Sigma^* \) be an acyclic morphism. Then \( \Delta^+_\mu(w) \geq 0 \) for every \( w \in \{a,b\}^* \).

Proof. Let \( \mu \equiv (\alpha, \beta) \). Since \( r_{bwt}(w) = r_{bwt}(w^m) \) for every \( w \in \Sigma^* \) and \( m > 1 \), let us assume that \( w \) is primitive. For the proof, we assume that \( |\alpha| \geq |\beta| \), and the other case is treated symmetrically. First, let us consider the case where \( \beta \) is not a suffix of \( \alpha \). Let moreover \( x \in \Sigma^* \) be the longest common suffix between \( \alpha \) and \( \beta \). It follows that there exist \( \alpha', \beta' \in \Sigma^* \) such that \( \alpha = \alpha'x \) and \( \beta = \beta'x \), and that the last symbol of \( \alpha' \) is different from the last of \( \beta' \) (otherwise \( x \) would be longer). Let \( R_x(\mu(w)) \) denote the multiset of rotations of \( \mu(w) \) with \( x \) as a prefix. Note that if \( x = \varepsilon \), then \( R_x(\mu(w)) = I_{\mu}(w) \). Since \( x \) appears in both \( \alpha \) and \( \beta \), it follows that \( |R_x(\mu(w))| \geq |w| \). Specifically, for each \( i \in [1,|w|] \), there exists \( t_i \in R_x(\mu(w)) \) such that \( t_i = x\mu(w[i+1,|w|] \cdot w[1,i-1]v) \), where \( v \) is either \( \alpha' \) or \( \beta' \), depending on whether \( w[i] \) is \( a \) or \( b \) respectively. The lexicographical order of these \( |w| \) rotations of \( \mu(w) \) with the same prefix correspond to the lexicographical order of the rotations in \( I_{\mu}(w) \), since by Corollary 8 the words \( \bigcup_{i=1}^{|w|} \{\mu(w[i+1,|w|] \cdot w[1,i-1])\} \) must differ in at least one position. By Corollary 12 this is either in the same or in the reverse order with respect to the sorting of the rotations of \( w \). Thus, there exists an injective coding \( \lambda : \{a,b\}^* \rightarrow \Sigma^* \subseteq \Sigma \) such that either \( \lambda(bwt(\mu(w))) \) or \( \lambda(bwt(\mu(w))^R) \) is a subsequence of \( bwt(\mu(w)) \), and therefore \( \Delta^+\mu(w) \geq r_{bwt}(w) \).

Let us now consider the case where \( \beta \) is suffix of \( \alpha \). Then, there exists a primitive word \( u \in \Sigma^+ \) and two integers \( p \geq q \geq 1 \) such that \( \beta = u^q \) and \( \alpha = \alpha'u^p \), with \( \alpha' \in \Sigma^+ \) that does not have \( u \) as suffix. Note that \( \alpha' \neq \varepsilon \), otherwise we would have \( \alpha = u^p \alpha' = u^p u^q = \beta \alpha \), i.e. \( \mu \) would not be acyclic. Let \( x \) be the longest common suffix between \( \alpha' \) and \( u \). If \( x \neq \alpha' \), from analogous arguments to the case where \( \beta \) is not a suffix of \( \alpha \), we have at least \( r_{bwt}(w) \) equal-letter runs in \( R_{xyv}(\mu(w)) \). Otherwise, if \( x = \alpha' \), let us consider the word \( y \in \Sigma^+ \) such that \( u = yx \). We can then consider the longest common suffix \( x' \) between \( xy \) and \( yx \), which must be a proper suffix (otherwise \( u \) would not be primitive), and apply the same reasoning over the set \( R_{x'yv}(\mu(w)) \) and the thesis follows.

The following example shows that Theorem 14 does not hold in the case of larger alphabets.

Example 15. Consider the acyclic morphism \( \mu \equiv (b,a,c) \). Then, \( bwt(beba) = beab \) and \( bwt(\mu(beba)) = bwt(acab) = cbaa \).

An immediate consequence of Theorem 14 is the following.
Corollary 16. Let $\mu : \{a, b\}^* \mapsto \Sigma^*$ be an acyclic morphism. Then, $\Delta_{\mu}^+(w) \geq 1$, for every $w \in \{a, b\}^*$.

The following theorem provides a characterization of cyclic morphisms in terms of the number of BWT-runs.

Theorem 17. A morphism $\mu : \{a, b\}^* \mapsto \Sigma^*$ is cyclic if and only if there exists $k > 0$ such that $r_{\text{bwt}}(\mu(w)) = k$ for all $w \in \{a, b\}^*$.

Proof. If $\mu \equiv (\alpha, \beta)$ is cyclic then there exists a primitive word $u \in \Sigma^*$ such that $\alpha = u^p$ and $\beta = u^q$, for some $p, q \geq 0$. Therefore, for each word $w \in \{a, b\}^*$, we have $r_{\text{bwt}}(\mu(w)) = r_{\text{bwt}}(u^p|w|_a + q|w|_b) = r_{\text{bwt}}(u)$. The other implication is a consequence of Theorem 14. In fact, by contraposition for each $k > 0$ we can find a word $w$ such that $r_{\text{bwt}}(w) > k$ (for instance, the $i$-th Thue–Morse finite word such that $i > \frac{k}{2}$ [4]), which leads to $r_{\text{bwt}}(\mu(w)) \geq r_{\text{bwt}}(w) > k$ as well.

4 Binary morphisms preserving $r_{\text{bwt}}$

This section is devoted to characterizing binary morphisms such that the number of BWT equal-letter runs is preserved after the action of the morphism on any binary word. First, we show with an example that this property is not trivial.

Example 18. Let $\theta \equiv (ab, aa)$ be the period-doubling morphism. It can be verified that $\Delta_{\theta}^+(ab) = 0$, $\Delta_{\theta}^+(aab) = 2$, and $\Delta_{\theta}^+(aaabaabab) = 4$.

Next, we show that every Sturmian morphism fixes the number of BWT-runs. From the definition of $E$, $\varphi$, and $\tilde{\varphi}$, and by Lemma 11, we derive the following.

Lemma 19. Let $w \in \{a, b\}^*$ be a binary word. Then, for all pairs of rotations $u$ and $v$ of $w$, and for each $\chi \in \{E, \varphi, \tilde{\varphi}\}$, it holds that $u < v$ if and only if $\chi(u) > \chi(v)$.

We prove that the number of BWT-runs is preserved by the morphisms that are the generators of the Sturmian morphisms. Note that from the following lemma a method can be derived to construct $\text{bwt}(\mu(w))$ starting from $\text{bwt}(w)$, for every Sturmian morphism $\mu$ and every binary word $w$.

Lemma 20. Let $w \in \{a, b\}^*$ be a binary word with $|\text{alph}(w)| = 2$. Then, for all $\chi \in \{E, \varphi, \tilde{\varphi}\}$, one has $r_{\text{bwt}}(w) = r_{\text{bwt}}(\chi(w))$. More in detail, one has $\text{bwt}(E(w)) = \text{bwt}(w)^R$ and $\text{bwt}(\varphi(w)) = \text{bwt}(\tilde{\varphi}(w)) = \text{bwt}(w)^R \cdot a^{|w|}$.

Proof. Since for each word $w$ and each integer $k > 0$ we have $r_{\text{bwt}}(w) = r_{\text{bwt}}(w^k)$, let us assume that $w$ is a primitive word. From Lemma 19, the case $\chi = E$ is trivial: in fact, from it follows that $\text{bwt}(E(w)) = \text{bwt}(w)^R$, and therefore $r_{\text{bwt}}(w) = r_{\text{bwt}}(E(w))$.

For the case $\chi = \varphi$ one can observe that every $b$ that occurs in $\varphi(w)$ is obtained from $\varphi(a)$, and therefore it is always preceded by an $a$. Thus, the rotations of $\varphi(w)$ left to cover are all those starting with an $a$, which therefore must also start with either $\varphi(a)$ or $\varphi(b)$. By Lemma 19, and by observing that $\varphi(a)$ ends with a $b$ and $\varphi(b)$ ends with $a$, we have that $\text{bwt}(\varphi(w)) = \text{bwt}(w)^R \cdot a^{|w|}$. Thus, we need to check if the run of $a$’s at the end merges with the last symbol of $\text{bwt}(w)^R$. This is equivalent to checking that the first symbol of $\text{bwt}(w)$ is a $b$, and by contradiction if the first rotation in lexicographical order is $ua$ for some $u \in \{a, b\}^{n-1}$, then $ua$ is a conjugate of $w$ and $ua < ua$ for each binary word $w$, a contradiction.
The next step after characterizing Sturmian morphisms as those fixing BWT equal-letter runs on binary words, is to find other binary morphisms that increase the number of BWT-runs always by the same fixed constant. Remind that if such a constant exists, it has to be an even integer because the BWT of any binary word starts with \( b \) and ends with \( a \).

We show that for every \( k > 0 \), we can find a morphism increasing the BWT-runs of any binary word by exactly \( 2k \). We do so by showing a family of binary morphisms that increase the BWT-runs always by \( 2 \), which then we can compose as we want. This family is formed by binary morphisms that are similar to the famous Thue–Morse morphism \( \tau \equiv (ab, ba) \).

\[ \begin{array}{c|c|c|c}
  M(w) & M(\varphi(w)) & M(\tilde{\varphi}(w)) \\
  \hline
  aabb\ b & a.a.ab.a.a.a & a.a.ba.a.ba.b & b \\
  ab\ a & a.a.b.a & a.a.ba.b & b \\
  ab\ a & a.a.b.a.a.a & a.a.ba.b.a.a & b \\
  ba\ a & a.a.a.a.a & a.a.ba.a.a.a & b \\
  ba\ a & a.a.a.a.a & a.a.ba.a.a.a & b \\
  b\ a & a.a.a.a.a & a.a.ba.a.a.a & b \\
  b\ a & a.a.a.a.a & a.a.ba.a.a.a & b \\
  b\ a & a.a.a.a.a & a.a.ba.a.a.a & b \\
  \end{array} \]

\[ \text{Figure 1 From left to right, the BWT-matrix for the words } w = \text{aabb}\ b, \varphi(w), \text{and } \tilde{\varphi}(w) \text{ respectively. For } M(\varphi(w)) \text{ and } M(\tilde{\varphi}(w)), \text{we separate with dots the images of symbols from } w. \text{ The rotations in bold of } M(\varphi(w)) \text{ and } M(\tilde{\varphi}(w)) \text{ correspond to the words in } I_\varphi(w) \text{ and } I_{\tilde{\varphi}}(w) \text{ respectively. The block of rotations in gray at the end of both } M(\varphi(w)) \text{ and } M(\tilde{\varphi}(w)) \text{ are in correspondence of the equal-letter run of } a \text{'s of length } |w|_a, \text{ which occurs for every } w \in \{a, b\}^* \text{. One can see that } \text{bwt}(\varphi(w)) = \text{bwt}(\tilde{\varphi}(w)) = \text{bwt}(w)^R \cdot a^{|w|_a}. \]

The following theorem shows a new characterization of Sturmian morphisms.

\textbf{Theorem 21.} Let \( \mu \) be a binary morphism. Then, the following are equivalent:
1. \( \Delta^I \mu(w) = 0 \) for every word \( w \) with \(|\text{alph}(w)| = 2\);
2. \( \mu \) is a Sturmian morphism.

\textbf{Proof.} By Theorem 3 and Lemma 20, all Sturmian morphisms preserve the number of BWT-runs. Conversely, suppose that \( \mu \) preserves the number of BWT-runs. By Theorem 17, such a morphism must be acyclic. Let \( s = \lim s_i \) be a characteristic Sturmian word. For every \( i \), the word \( \mu(s_i) \) has 2 runs in its BWT, hence it is circularly balanced (Proposition 4). Let us consider the word \( \mu(s) = \lim \mu(s_i) \). It is balanced and aperiodic, since it is obtained by applying an acyclic morphism to a Sturmian word \([6]\). Then, \( \mu(s) \) is Sturmian by using Theorem 1, whence \( \mu \) is a Sturmian morphism by applying Theorem 2.

\section{Binary morphisms increasing \( r_{\text{bwt}} \) by a constant}

The next step after characterizing Sturmian morphisms as those fixing BWT equal-letter runs on binary words, is to find other binary morphisms that increase the number of BWT-runs always by the same fixed constant. Remind that if such a constant exists, it has to be an even integer because the BWT of any binary word starts with \( b \) and ends with \( a \).

We show that for every \( k > 0 \), we can find a morphism increasing the BWT-runs of any binary word by exactly \( 2k \). We do so by showing a family of binary morphisms that increase the BWT-runs always by \( 2 \), which then we can compose as we want. This family is formed by binary morphisms that are similar to the famous Thue–Morse morphism \( \tau \equiv (ab, ba) \). The
structure of the BWT of Thue–Morse words has been studied before and it is well understood [4, 10]. We generalize such results by showing how to derive $\text{bwt}(\mu(w))$ from $\text{bwt}(w)$ for every Thue–Morse-like morphism $\mu$ and every binary word $w$.

**Definition 22.** A binary morphism is Thue–Morse-like if it has the form $\tau_{p,q} \equiv (ab^p, ba^q)$ for some $p, q > 0$.

We prove the following proposition, which is crucial to obtain the main result of this section. Figure 2 highlights the key aspects of the proof.

**Proposition 23.** For every binary word $w$ such that $\text{alph}(w) = \{a, b\}$, the I-rotations of $\tau_{p,q}(w)$ are contiguous in the BWT-matrix of $\tau_{p,q}(w)$, and their last letters spell $\text{bwt}(w)$.

**Proof.** Let $w$ be a binary word of length $n$ such that $\text{alph}(w) = \{a, b\}$. Observe that $\tau_{p,q} \equiv (ab^p, ba^q)$ is abelian order-preserving, so the I-rotations of $\tau_{p,q}(w)$ maintain their relative order. Because $\tau_{p,q}(a)$ ends with $b$ and $\tau_{p,q}(b)$ ends with $a$, if we consider only the I-rotations of $\tau_{p,q}(w)$ and take the last letter of each, we obtain $\text{bwt}(w)$, which starts with $a$ and ends with $b$. It remains to show that all the I-rotations of $\tau_{p,q}(w)$ are contiguous in its BWT-matrix.

If $p > 1$, each S-rotation starting with $a$, has to start either with $ab$ for some $2 \leq i \leq q+1$, or with $aba^q$, and both of these prefixes are smaller than $ab^p$. If $p = 1$, an S-rotation starting with $a$ is smaller than the word $a(ba^q)^{n-1}ba^{q-1}$, which is smaller than the rotation having $(ab)^q ba$ as a prefix for some $0 < i < n$. The I-rotations that start with $a$ have prefixes of such type. In both cases, we obtain that the S-rotations starting with $a$ are smaller than the I-rotations starting with $a$. A symmetric argument shows that S-rotations starting with $b$ are greater than the I-rotations starting with $b$. Thus, the I-rotations are contiguous and the thesis holds.

Now we are ready to show that Thue–Morse-like morphisms increase the number of BWT-runs of binary words always by 2.

**Lemma 24.** For every binary word $w$ such that $\text{alph}(w) = \{a, b\}$, it holds that

$$\text{bwt}(\tau_{p,q}(w)) = b^{\text{len}(w)-1} \cdot \text{bwt}(w) \cdot b^{(p-1)\text{len}(w)} a^{\text{len}(w)}$$

and that $r_{\text{bwt}}(\tau_{p,q}(w)) = r_{\text{bwt}}(w) + 2$.

**Proof.** We show that the block of $\text{bwt}(\tau_{p,q}(w))$ that corresponds to the S-rotations starting with the letter $a$ is equal to $b^{\text{len}(w)-1}a$. If $q = 1$, all the S-rotations starting with $a$ end with the letter $b$. If $q > 1$, the only S-rotations that start with $a$ and end with $b$ have as a prefix either $a^{q-1}b$ or $a^q ba$. The smallest S-rotation starting with $a$ and ending with $a$ starts with $a^q ba$. Hence, S-rotations starting with $a$ and ending with $b$ appear before those ending with $a$.

It follows that the block of $\text{bwt}(\tau_{p,q}(w))$ defined by the S-rotations starting with $a$ spells $b^{\text{len}(w)-1}a$, because of their order, and because each of these rotations is in correspondence with some specific $a$ inside $\tau_{p,q}(b)$ for some specific $b$ of $w$. Only one of these $a$’s per image produces a rotation ending with $b$, and the other $q - 1$ $a$’s yield rotations ending with $a$.

Showing that the block of $\text{bwt}(\tau_{p,q}(w))$ corresponding to the S-rotations starting with the letter $b$ equals $b^{(p-1)\text{len}(w)} a^{\text{len}(w)}$ is handled symmetrically.

By using Proposition 23, we obtain

$$\text{bwt}(\tau_{p,q}(w)) = b^{\text{len}(w)-1} \cdot \text{bwt}(w) \cdot b^{(p-1)\text{len}(w)} a^{\text{len}(w)}$$

As $\text{bwt}(w)$ starts with $a$ and ends with $b$, we have that $r_{\text{bwt}}(\tau_{p,q}(w)) = r_{\text{bwt}}(w) + 2$, and the thesis holds.
As a consequence of Theorem 21 and Lemma 24, we obtain the following corollary.

\begin{corollary}
Given a non-negative even integer \(2t\), there exists a binary morphism \(\mu\) such that \(\Delta^+ \mu (w) = 2t\) and \(\Delta^\times \mu (w) \leq t + 1\), for every word \(w\) with \(|\text{alph}(w)| = 2\).
\end{corollary}

\textbf{Proof.} We can construct the morphism \(\mu \in (\{E, \varphi, \tilde{\varphi}\} \cup \{(ab^p, ba^q) | p, q > 0\})^*\) such that \(\mu\) is obtained by composing, in any order, exactly \(t\) morphisms taken in the set \(\{(ab^p, ba^q) | p, q > 0\}\) and an arbitrary number of Sturmian morphisms. By Theorem 21 and Lemma 24, it holds that \(\Delta^+ \mu (w) = 2t\). The value of the function \(\Delta^\times \mu (w) = (r_{bwt}(w) + 2t)/r_{bwt}(w) = 1 + 2t/r_{bwt}(w)\) is maximized when \(r_{bwt}(w) = 2\). This maximum is \(\Delta^\times \mu (w) = t + 1\).

We conclude this section by showing a simple algorithm that allows us to construct an arbitrarily large family of words \(w_1, w_2, \ldots\) with exactly \(2t\) BWT-runs each. In Algorithm 1, a morphism \(\mu\) such that \(\Delta^+ \mu (w) = 2(t - 1)\) for every binary word is required. Note that Corollary 25 assures that such a morphism exists.

Moreover, each word \(w_i\) is a prefix of the next word \(w_{i+1}\), so that the infinite word \(w = \lim_{i \to \infty} w_i\) is well defined, and it is aperiodic. This is given because it holds for the (implicit) standard Sturmian words \(s_i\) for \(i \in [1, k]\) being used, which are circularly balanced (i.e., \(r_{bwt}(w) = 2\) on them, as reported in Proposition 4), and their limit is a characteristic Sturmian word, which is aperiodic.
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![Algorithm 1 Algorithm for constructing words with 2t BWT-runs.](image)

**Require:** A morphism $\mu$ with $\Delta_\mu^+(w) = 2(t - 1)$. A sequence of positive integers $d_1, \ldots, d_k$.

**Ensure:** A sequence of words $w_1, w_2, \ldots, w_k$ where $r_{\text{bwt}}(w_i) = 2t$ for any $1 \leq i \leq k$.

\[
\begin{align*}
  w_{i-1} & \leftarrow \mu(b) \\
  w_0 & \leftarrow \mu(a) \\
  \text{for } i \in [1, k] & \text{ do} \\
  & w_i \leftarrow w_{i-1}w_{i-2} \\
  \text{end for} \\
  \text{return } w_1, \ldots, w_k
\end{align*}
\]

6 Morphisms with an unbounded increase on $r_{\text{bwt}}$

There exist morphisms that do not behave as well as Sturmian and Thue–Morse-like morphisms with respect to $r_{\text{bwt}}$. If we consider an alphabet of size greater than 2, we can always find a morphism $\mu$ such that the values $\Delta_\mu^+(w)$ and $\Delta_\mu^+(w)$ are arbitrarily large.

► **Lemma 26.** Let $\Sigma = \{c_1, \ldots, c_k, a, b\}$ with $k \geq 1$. Let $\varphi \equiv (ab, a)$ be the Fibonacci morphism. Then, $r_{\text{bwt}}(w) = k + 3$ if $w$ belongs to $\{\varphi^{2i}(a)c_1c_2\cdots c_k[i \geq 1]\}$.

**Proof.** We prove the result by induction on $k \geq 1$. Observe that the words $\varphi^{2i}(a)$ for $i \geq 1$ are Fibonacci words ending with the letter $a$. It is known that in these words, if we append the letter $c_1$ smaller than $a$ at the end, then the number of runs becomes 4 [31, Theorem 11]. For the inductive step, suppose that $r_{\text{bwt}}(\varphi^{2i}(a)c_1 \cdots c_{k-1}) = k + 2$. When appending $c_k$ at the end, the rotations that do not start with $c_k$ keep their relative order, and the rotation that originally ended with $c_{k-1}$ now ends with $c_k$. Hence, they define the same number of runs as before. The rotation starting with $c_k$ can be found after the rotation starting with $c_{k-1}$, which does not end with $b$, and before the first rotation starting with $a$, which ends with $b$. Hence, the number of runs increases by 1. Thus, $r_{\text{bwt}}(\varphi^{2i}(a)c_1 \cdots c_k) = k + 3$.

► **Lemma 27.** Let $\Sigma = \{c_1, \ldots, c_k, a, b\}$ with $k \geq 1$. Let $\varphi \equiv (ab, a)$ be the Fibonacci morphism, and $\mu \equiv (c_1, c_2, \ldots, c_k, ab, a)$ be a morphism on the alphabet $\Sigma$. Then, $r_{\text{bwt}}(w) = \Omega(\log n)$ for every $w \in \{\mu(\varphi^{2i}(a)c_1c_2\cdots c_k[i \geq 1]\}$.

**Proof.** The morphism $\mu$ maps a Fibonacci word ending with $a$ having $c_1 \ldots c_k$ appended at the end, to the next Fibonacci word, which ends with $b$, having $c_1 \ldots c_k$ appended at the end. For $k = 1$, it is known that the number of runs in this family is $\Omega(\log n)$ [15]. In a similar way to Lemma 26, it is possible to prove by induction that appending $c_k$ at the end of $\mu(\varphi^{2i}(a)c_1c_2\cdots c_{k-1})$ adds 2 runs when $k = 2$ and exactly 1 new run when $k > 2$.

► **Proposition 28.** For each alphabet $\Sigma$ with size greater than 2 there exist a morphism $\mu$, satisfying that for every $k$, there is a word $w \in \Sigma^*$ such that $\Delta_\mu^+(w) \geq k$ and $\Delta_\mu^+(w) \geq k$.

**Proof.** This is immediate from Lemma 26 together with Lemma 27.

Finding examples like the previous ones for binary morphisms is trickier, but at least in the case of $\Delta_\mu^+$, it is possible. An example of a binary morphism for which the value $\Delta_\mu^+(w)$ can be arbitrarily large is the period-doubling morphism denoted by $\theta$ and defined by the rules $\theta(a) = ab$ and $\theta(b) = aa$.

► **Lemma 29.** Let $\theta$ be the period-doubling morphism. For any positive integer $k$ there exist a word $w$ such that $\Delta_\theta^+(w) > k$. 
Proof. W.l.o.g assume that $k > 2$. For $i \in [2, k]$ define the words
$s_i = ab^i a \cdot u_i$ and $e_i = ab^i a \cdot u_i^R$, where $u_i = a^{2k-i}ba^{i-2}$.

We say that $s_i$ is a starting factor, and $e_i$ is an ending factor. Observe that $s_i$ (resp. $u_i$) is always smaller than $e_i$ (resp. $u_i^R$). Moreover, it holds that if $i < j$, then $u_i < u_j < u_j^R < u_i^R$.

We define the word $w_k = (\Pi_{i=2}^k s_i) a^k$ and show that $\Delta_p^+(w_k) = 2k$. Figure 3 shows the structure of both BWTs and highlights the increase in the number of runs.

\begin{center}
\begin{tabular}{|c|c|c|c|}
\hline
$a$ & $...$ & $...$ & $b$
\hline
$...$ & $...$ & $x$
\hline
$a$ & $b^k a u_k^R$ & $...$ & $a$
\hline
$ba$ & $a^{2k} b$ & $...$ & $a$
\hline
$...$ & $...$ & $y$
\hline
$ba$ & $bb$ & $...$ & $a$
\hline
$b^2 a$ & $u_2$ & $...$ & $a$
\hline
$b^2 a$ & $u_3$ & $...$ & $ab$
\hline
$...$ & $...$ & $...$
\hline
$b^2 a$ & $u_k$ & $...$ & $ab^{k-2}$
\hline
$b^2 a$ & $u_k^R$ & $...$ & $ab^{k-2}$
\hline
$...$ & $...$ & $...$
\hline
$b^2 a$ & $u_2^R$ & $...$ & $ab$
\hline
$b^2 a$ & $u_2^R$ & $...$ & $a$
\hline
$...$ & $...$ & $...$
\hline
$b^2 a$ & $u_k$ & $...$ & $a$
\hline
$b^2 a$ & $u_k^R$ & $...$ & $a$
\hline
\end{tabular}
\end{center}

\[\theta^{\equiv(ab, a)}\]

**Figure 3** To the left is the BWT-matrix of $w_k$. To the right is the BWT-matrix of $\theta(w_k)$, here displayed in reverse order. Each row represents a block of rotations from $S_0(w_k)$ starting with the same prefix, highlighted in the first column. Each one of these block except the first one yields 2 extra runs on $\text{bwt}(\theta(w_k))$. The words $x$ and $y$ correspond to the concatenation of the last letters of blocks of the BWT-matrix of $w_k$ whose form is unknown, but do not play a role in the increase on $\Delta_p^+(\theta(w_k))$.

Consider the rotations of $w_k$ starting with $b^i a$ with $1 < i \leq k$. The left shift of the unique rotation starting with the $i$-th starting factor, and the left shift of the unique rotation starting with the $i$-th ending factor, are the smallest and greater, respectively. Both of them end with the letter $a$. The remaining rotations starting with $b^i a$ (if any) have to end with $b$ because in them the prefix $b^i a$ corresponds to a suffix of a longer run of $b$'s followed by an $a$.

In the case of the rotations of $w_k$ starting with $ba$, the one starting in the last $b$ of $e_k$, has $ba^k a^k$ as a prefix, so it is the smallest of them. Also, this rotation is preceded by the factor $ab^k a u_k^{b-2}$, which ends in $a$. The greatest rotation starting with $ba$ is the one starting with the $b$ preceding $e_2$, which is followed by $abb$ and preceded also by an $a$. In the case of the rotations of $w_k$ starting with $a$, the smallest of them ends with the letter $b$ as in any binary word. The greatest is the rotation starting with $ab^k a u_k^R$ which is preceded by the letter $a$. 

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With the general structure of the BWT of $w_k$ in mind, now we analyse the BWT of $\theta(w_k)$. The morphism $\theta$ is order-reversing and all the I-rotations of $\theta(w_k)$ start with the letter $a$. S-rotations of $\theta(w_k)$ starting with the letter $b$ are always preceded by an $a$, and it is easy to see that this run of $a's$ merges with the last $a$ in the greatest I-rotation. The S-rotations starting with an $a$ have an even number of $a's$ before the first $b$ appears, and also end with the letter $a$. This implies that they appear grouped after all the I-rotations of the form $(aa)^i ab$ for some $1 \leq i \leq k$, and before all the I-rotations starting with $(aa)^{i-1} ab$. As the smallest and greatest rotations of each of these blocks of I-rotations end with $b$ (because of the action of $\theta$), it follows that the group of S-rotations starting with $(aa)^i b$ increases the number of runs of the BWT of $\theta(w_k)$ by 2 with respect to the BWT of $w_k$. This happens for $1 \leq i \leq k$, so the overall increase in $r_{\text{bwt}}$ after applying the morphism $\theta$ is exactly $2k$. ▶

From the lemma above we can deduce that there are binary morphisms that can greatly increase the number of BWT-runs of some words. We define the sensitivity of BWT-runs to morphism application in a similar way to how Akagi et al. define the sensitivity of repetitiveness measures to edit operations [1].

Definition 30. The BWT additive sensitivity and BWT multiplicative sensitivity for a morphism $\mu$ are respectively, the functions

$$A_{\mu}(n) = \max_{w \in \Sigma^n} (\Delta_{\mu}^+(w)) \text{ and } M_{\mu}(n) = \max_{w \in \Sigma^n} (\Delta_{\mu}^x(w))$$

Proposition 31. Let $\theta$ be the period-doubling morphism. It holds that $A_{\theta}(n) = \Theta(\sqrt{n})$.

Proof. The length of the words $w_k$ in Lemma 29 is $n = \Theta(k^2)$. We showed that $\Delta_{\theta}^+(w_k) = 2k = \Theta(\sqrt{n})$ on these words. For values of $n$ in between $|w_k|$ and $|w_{k+1}|$, it is easy to see that for the word $w_k a^j$ for $0 < j < |w_{k+1}| - |w_k|$, it still holds that $\Delta_{\theta}^+(w_k a^j) = 2k$, as none of the key aspects of the proof of Lemma 29 changes. Thus, the claim is true. ▶

7 On the impact of morphisms on other repetitiveness measures

Morphisms behave very differently when other repetitiveness measures are considered. For a general survey on repetitiveness measures see [29]. For instance, any morphism $\mu$ increases the size of the Lempel-Ziv parsing [22] of any word by at most an additive constant depending only on $\mu$. This holds for any alphabet size, as shown by Constantinescu and Ilie [9, Lemma 8].

Lemma 32. Let $\mu : \Sigma^* \rightarrow \Gamma^*$ be any morphism. For every word $w$, it holds that $z(\mu(w)) \leq z(w) + k$ where $k$ is a constant depending only on $\mu$.

The result of Constantinescu and Ilie can easily be extended to the LZ parsing without overlaps [22], the optimal (not the greedy) LZ-end parsing [20], and bidirectional macro-schemes [38]. We can show a similar result for the measure $g(w)$ defined as the size of the smallest deterministic context-free grammar generating only $w$ [18]. This can be further generalized to the size of the smallest run-length context-free grammar [33], and also to the size of the smallest collage system [17].

Lemma 33. Let $\mu : \Sigma^* \rightarrow \Gamma^*$ be any (possible erasing) morphism. For every word $w$, it holds that $g(\mu(w)) \leq g(w) + k$ where $k$ is a constant depending only on $\mu$.

Proof. Given a deterministic context-free grammar $G$ of size $|G|$ generating $w$, we construct a grammar generating $\mu(w)$. For each occurrence of a terminal symbol $a$ in any rule of the grammar, replace it with a new non-terminal $A_a$. For each terminal symbol add the rule
Then, $\rho$ every word Thue–Morse-like and Sturmian morphisms are the only binary morphisms increasing which increase the number of BWT-runs of binary words by 2. The following question is left open.

Fact, we showed an infinite family of binary morphisms called Thue–Morse-like morphisms, us to construct binary words with any possible number of BWT-runs, and morphisms with its own, when this characterization is in conjunction with the rest of our results, it allows number of BWT-runs for all words equal-letter runs of finite words.

In this work, we have studied the impact of morphism application on the number of BWT runs of purely morphic words, this is known to hold only for the binary case [12].

In other variables by erasing) morphism has erasing rules, we can delete them, and replace the occurrences of those erasing variables in other variables by $\varepsilon$. We repeat this recursively. The size of the resulting grammar can only decrease, so the thesis still holds.

If for some fixed measure and morphism, this morphism increases the value of the measure always by at most a fixed constant, then we can derive an easy upper bound for the family of words obtained by iterating that morphism.

Proposition 34. Let $\rho$ be a repetitiveness measure and $\mu$ be a morphism. Suppose that for every word $w$ it holds that $\rho(\mu(w)) \leq \rho(w) + k$ for a constant $k$ depending only on $\rho$ and $\mu$. Then, $\rho = O(i)$ in the family $\{\mu^i(w) \mid i \geq 0\}$.

Proof. Let $k' = \rho(\mu(w))$. We show by induction that $\rho(\mu^i(w)) \leq ki + k'$ for any $i \geq 1$. For $i = 1$, clearly $\rho(\mu(w)) \leq k + k'$. Let $i > 1$ and suppose the claim is true for $i - 1$. Then, $\rho(\mu^i(w)) \leq \rho(\mu^{i-1}(w)) + k \leq (k(i - 1) + k') + k \leq ki + k'$.

The families on the proposition above are known as D0L-sequences [36]. As a direct consequence of Lemma 33 and Proposition 34, it holds that all repetitiveness measures upper-bounded by $g$ are $O(i)$ on the family of words belonging to a fixed D0L-sequence. In fact, the result we obtain is even more general because we can apply any morphism to words obtained from a D0L-sequence increasing the size of the grammar only by a fixed constant.

Proposition 35. For every (possibly erasing) morphisms $\mu$ and $\lambda$, and every word $w$, it holds that $g = O(i)$ in the family $\{\lambda \circ \mu^i(w) \mid i \geq 0\}$.

Proof. By Lemma 33 and Proposition 34, it holds that $g(\mu^i(w)) = O(i)$ for every (possibly erasing) morphism $\mu$. By Lemma 33, one has $g(\lambda \circ \mu^i(w)) \leq g(\mu^i(w)) + k$ for every (possibly erasing) morphism $\lambda$, and a constant $k$ depending on $\lambda$. Thus, $g(\lambda \circ \mu^i(w)) = O(i)$.

It is unknown if an analogous result is true for $r_{\text{bwt}}$. In fact, even for the restricted case of purely morphic words, this is known to hold only for the binary case [12].

8 Conclusions and further work

In this work, we have studied the impact of morphism application on the number of BWT equal-letter runs of finite words.

Firstly, we characterized Sturmian morphisms as the binary morphisms preserving the number of BWT-runs for all words $w$ such that $|\text{alph}(w)| = 2$. Besides being interesting on its own, when this characterization is in conjunction with the rest of our results, it allows us to construct binary words with any possible number of BWT-runs, and morphisms with known behavior. This can have practical applications, for instance, in experimentation. In fact, we showed an infinite family of binary morphisms called Thue–Morse-like morphisms, which increase the number of BWT-runs of binary words by 2. As a consequence, we have extended the results of Brlek et al. [4] on the number of BWT-runs of words generated by iterating the composition of the Fibonacci morphism with the Thue–Morse morphism to any composition of Sturmian morphisms and Thue–Morse-like morphisms. Also, we are able to construct infinite sequences of words of increasing length, having all exactly $2k$ BWT-runs, and converging to an aperiodic infinite word at their limit. While the result on Sturmian morphisms is a complete characterization, it is unknown if the compositions of Thue–Morse-like and Sturmian morphisms are the only binary morphisms increasing the number of BWT-runs exactly by 2. The following question is left open.
Question 36. What is a sufficient and necessary condition for a binary morphism $\mu$, to have $\Delta^\times_{\mu}(w) \leq 2k$ (where $k > 0$) for every binary word $w$?

We showed that when the alphabet size of the domain is $\sigma > 2$, the values $\Delta^+_{\mu}(w)$ and $\Delta^\times_{\mu}(w)$ can be arbitrarily large for some morphisms. In the case of the binary alphabet, we went further and showed that there exists morphisms where $AS_{\mu}(w) = \Omega(\sqrt{n})$. We plan to extend such a result by studying morphisms where all the images of letters are primitive words. On the other hand, it is unknown if the value $\Delta^\times_{\mu}(w)$ can be unbounded for some morphism $\mu$ with binary domain. We conjecture that this is not the case.

Conjecture 37. For every morphism $\mu : \{a,b\}^* \mapsto \Sigma^*$, we can find a constant $k$ such that $\Delta^\times_{\mu}(w) \leq k$, for every word $w \in \{a,b\}^*$.

If Conjecture 37 were true, the following conjecture on images of standard Sturmian words would also be true.

Conjecture 38. For every morphism $\mu : \{a,b\}^* \mapsto \Sigma^*$ and every sequence of standard Sturmian words $(s_i)_{i \in \mathbb{N}}$, it holds that $r_{bwt}(\mu(s_i)) = \Theta(1)$.

Finally, we showed that the impact of morphism application on BWT-runs, is quite different from the impact of morphisms on other repetitiveness measures based on popular compression schemes, like context-free grammars and LZ factorizations. In these measures, the additive increase after morphism application is bounded by a constant depending only on the morphism and the measure. This raises the following question, which is true in the case of smallest grammars and (some) variants of the LZ parsing, but unknown in the case of BWT equal-letter runs.

Question 39. Does it hold that $r_{bwt}(w) = O(i)$ when $w \in \{\lambda \circ \mu^i(a) \mid i \geq 0\}$, for every pair of morphisms $\mu : \{a,b\}^* \mapsto \{a,b\}^*$ and $\lambda : \{a,b\}^* \mapsto \Sigma^*$?

We are working on proving or refuting these questions and conjectures. In the future, we plan to study how to extend the results on morphism fixing BWT-runs, and morphisms increasing BWT-runs by a fixed natural number, to alphabets of size greater than 2.

References

10:18 On the Impact of Morphisms on BWT-Runs


Comparing Elastic-Degenerate Strings: Algorithms, Lower Bounds, and Applications

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Abstract
An elastic-degenerate (ED) string $T$ is a sequence of $n$ sets $T[1], \ldots, T[n]$ containing $m$ strings in total whose cumulative length is $N$. We call $n$, $m$, and $N$ the length, the cardinality and the size of $T$, respectively. The language of $T$ is defined as $L(T) = \{ S_i : S_i \in T[i] \text{ for all } i \in [1, n] \}$. ED strings have been introduced to represent a set of closely-related DNA sequences, also known as a pangenome. The basic question we investigate here is: Given two ED strings, how fast can we check whether the two languages they represent have a nonempty intersection? We call the underlying problem the ED STRING INTERSECTION (EDSI) problem. For two ED strings $T_1$ and $T_2$ of lengths $n_1$ and $n_2$, cardinalities $m_1$ and $m_2$, and sizes $N_1$ and $N_2$, respectively, we show the following:

- There is no $O((N_1 N_2)^{1-\epsilon})$-time algorithm, thus no $O((N_1 m_2 + N_2 m_1)^{1-\epsilon})$-time algorithm and no $O((N_1 N_2 + N_2 m_1)^{1-\epsilon})$-time algorithm, for any constant $\epsilon > 0$, for EDSI even when $T_1$ and $T_2$ are over a binary alphabet, unless the Strong Exponential-Time Hypothesis is false.
- There is no combinatorial $O((N_1 + N_2)^{1-2\epsilon} f(n_1, n_2))$-time algorithm, for any constant $\epsilon > 0$ and any function $f$, for EDSI even when $T_1$ and $T_2$ are over a binary alphabet, unless the Boolean Matrix Multiplication conjecture is false.
- An $O(N_1 \log N_1 \log n_1 + N_2 \log N_2 \log n_2)$-time algorithm for outputting a compact (RLE) representation of the intersection language of two unary ED strings. In the case when $T_1$ and $T_2$ are given in a compact representation, we show that the problem is NP-complete.
- An $O(N_1 m_2 + N_2 m_1)$-time algorithm for EDSI.
- An $O(N_1^{\omega-1} n_2 + N_2^{\omega-1} n_1)$-time algorithm for EDSI, where $\omega$ is the exponent of matrix multiplication; the $\tilde{O}$ notation suppresses factors that are polylogarithmic in the input size.

We also show that the techniques we develop have applications outside of ED string comparison.

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Comparing Elastic-Degenerate Strings: Algorithms, Lower Bounds, and Applications

1 Introduction

Sequence (or string) comparison is a fundamental task in computer science, with numerous applications in computational biology [24], signal processing [16], information retrieval [7], file comparison [25], pattern recognition [4], security [36], and elsewhere [37]. Given two or more sequences and a distance function, the task is to compare the sequences in order to infer or visualize their (dis)similarities [15].

Many sequence representations have been introduced over the years to account for unknown or uncertain letters, a phenomenon that often occurs in data that comes from experiments [8]. In the context of computational biology, for example, the IUPAC notation [27] is used to represent locations of a DNA sequence for which several alternative nucleotides are possible. This gives rise to the notion of degenerate string (or indeterminate string): a sequence of finite sets of letters [2]. When all sets are of size 1, we are in the special case of a standard string (or deterministic string). Degenerate strings can encode the consensus of a population of DNA sequences [17] in a gapless multiple sequence alignment (MSA). Iliopoulos et al. generalized this notion to also encode insertions and deletions (gaps) occurring in MSAs by introducing the notion of elastic-degenerate string: a sequence of finite sets of strings [26].

The main motivation to consider elastic-degenerate (ED) strings is that they can be used to represent a pangenome: a collection of closely-related genomic sequences that are meant to be analyzed together [42]. Several other, more powerful, pangenome representations have been proposed in the literature, mostly graph-based ones; see the comprehensive survey by Carletti et al. [12] or by Baaijens et al. [5]. Compared to these more powerful representations, ED strings have at least two algorithmic advantages, as they support: (i) fast and simple on-line string matching [23, 13]; and (ii) (deterministic) subquadratic string matching [3, 9, 10].

Our main goal here is to give the first algorithms and lower bounds for comparing two pangenomes represented by two ED strings.\(^1\) We consider the most basic notion of matching, namely, to decide whether two ED strings, each encoding a language, have a nonempty intersection. Like with standard strings, algorithms for pairwise ED string comparison can serve as computational primitives for many analysis tasks (e.g., phylogeny reconstruction); lower bounds for pairwise ED string comparison can serve as meaningful lower bounds for more powerful pangenome representations such as, for instance, variation graphs [12].

Let us start with some basic definitions and notation. An alphabet \(\Sigma\) is a finite nonempty set of elements called letters. By \(\Sigma^*\) we denote the set of all strings over \(\Sigma\) including the empty string \(\varepsilon\) of length 0. For a string \(S = S[1] \cdots S[n]\) over \(\Sigma\), we call \(n = |S|\) its length. The fragment \(S[i..j]\) of \(S\) is an occurrence of the underlying substring \(P = S[i] \cdots S[j]\). We also say that \(P\) occurs at position \(i\) in \(S\). A prefix of \(S\) is a fragment of \(S\) of the form \(S[1..j]\) and a suffix of \(S\) is a fragment of \(S\) of the form \(S[i..n]\). An elastic-degenerate string (ED string, in short) \(T\) is a sequence \(T = T[1] \cdots T[n]\) of \(n\) finite sets, where \(T[i]\) is a subset of \(\Sigma^*\). The total size of \(T\) is defined as \(N = N_\varepsilon + \sum_{i=1}^n |T[i]|\), where \(N_\varepsilon\) is the total number of empty strings in \(T\). By \(m\) we denote the total number of strings in all \(T[i]\), i.e., \(m = \sum_{i=1}^n |T[i]|\). We say that \(T\) has length \(n = |T|\), cardinality \(m\) and size \(N = ||T||\). An ED string \(T\) can be treated as a compacted nondeterministic finite automaton (NFA) with \(n+1\) states, numbered 1, \ldots, \(n+1\), and \(m\) transitions labeled by strings in \(\Sigma^*\). State 1 is starting and state \(n+1\) is accepting. For each index \(i \in [1,n]\) and string \(S \in T[i]\), there is a

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\(^1\) Pangenome comparison is one of the central goals of two large EU funded projects on computational pangenomics: PANGAIA (https://www.pangenome.eu/) and ALPACA (https://alpaca-itn.eu/).
transition from state $i$ to state $i+1$ with label $S$; inspect also Figure 1 for an example. The language $\mathcal{L}(T)$ generated by the ED string $T$ is the language accepted by this compacted NFA. That is, $\mathcal{L}(T) = \{S_1 \cdots S_n : S_i \in T[i] \text{ for all } i \in [1,n]\}$.

**MSA**

\[
\text{GTTCAGTTTAC--AA}
\]

\[
\text{GTTCAGTTTTACACAA}
\]

\[
\text{GTTCAGATT----AA}
\]

**ED string**

\[
\{\text{GTT}\}, \{\text{T}\}, \{\text{AG}\}, \{\text{TT}\}, \{\text{ACAC}\}, \{\text{AA}\}
\]

\[
\text{NFA}
\]

\[
\begin{array}{c}
\text{starting} \\
\text{AG} \\
\text{T} \\
\text{TT} \\
\text{ACAC} \\
\text{accepting}
\end{array}
\]

**Figure 1** An example of an MSA and its corresponding (non-unique) ED string $T$ of length $n = 7$, cardinality $m = 11$ and size $N = 20$, and the compacted NFA for $T$. The compacted NFA can be seen as a special case of an edge-labeled directed acyclic graph.

We next define the main problem in scope; inspect also Figure 2 for an example.

**ED String Intersection (EDSI)**

**Input**: Two ED strings, $T_1$ of length $n_1$, cardinality $m_1$ and size $N_1$, and $T_2$ of length $n_2$, cardinality $m_2$ and size $N_2$.

**Output**: YES if $\mathcal{L}(T_1)$ and $\mathcal{L}(T_2)$ have a nonempty intersection, NO otherwise.

**Figure 2** An example of two ED strings $T_1$ and $T_2$ with their parameters and the intersection of their languages. In this instance, we see that $\mathcal{L}(T_1)$ and $\mathcal{L}(T_2)$ have a nonempty intersection.

**Our Results.** We assume that the ED strings are over an integer alphabet $[1, (N_1 + N_2)^\mathcal{O}(1)]$. We make the following specific contributions:

1. In Section 2.1, we give several conditional lower bounds. In particular, we show that there is no $\mathcal{O}((N_1 N_2)^{1-\epsilon})$-time algorithm, thus no $\mathcal{O}((N_1 m_2 + N_2 m_1)^{1-\epsilon})$-time algorithm and no $\mathcal{O}((N_1 n_2 + N_2 n_1)^{1-\epsilon})$-time algorithm, for any constant $\epsilon > 0$, for EDSI even when $T_1$ and $T_2$ are over a binary alphabet, unless the Strong Exponential-Time Hypothesis (SETH) [11] is false.

2. In Section 2.2, we present other conditional lower bounds. In particular, we show that there is no combinatorial $\mathcal{O}((N_1 + N_2)^{1.2-\epsilon} f(n_1, n_2))$-time algorithm, for any constant $\epsilon > 0$ and any function $f$, for EDSI even when $T_1$ and $T_2$ are over a binary alphabet, unless the Boolean Matrix Multiplication (BMM) conjecture [1] is false.
3. In Section 3, we show an $O(N_1 \log N_1 \log n_1 + N_2 \log N_2 \log n_2)$-time algorithm for outputting a compact (RLE) representation of the intersection language of two unary ED strings. In the case when $T_1$ and $T_2$ are given in a compact representation, we show that the problem is NP-complete.

4. In Section 4.1, we show an $O(N_1 m_2 + N_2 m_1)$-time algorithm for EDSI.

5. In Section 4.2, we show an $\tilde{O}(N_1 \omega^{-1} n_2 + N_2 \omega^{-1} n_1)$-time algorithm for EDSI, where $\omega$ is the matrix multiplication exponent.

Interestingly, we show that the techniques we develop here have applications outside of ED string comparison. Given a sequence $P = P_1, \ldots, P_n$ of $n$ standard strings, we define an acronym of $P$ as a string $A = A_1 \cdots A_n$, where $A_i$ is a possibly empty prefix of $P_i$, for all $i \in [1, n]$. In the Acronym Generation (AG) problem, we are given a set $D$ of $k$ strings of total length $K$ and a sequence $P$ of $n$ strings of total length $N$, and we are asked to say YES if and only if some acronym of $P$ belongs to $D$. In Section 5, we show how our techniques for EDSI can be modified to solve AG in $O(nK + N)$ time.

Related Work. Apart from its applications to pangenome comparison, EDSI is interesting theoretically on its own as a special case of regular expression (regex) matching. Regex is a basic notion in formal languages and automata theory. Regular expressions are commonly used in practical applications to define search patterns. Regex matching and membership testing are widely used as computational primitives in programming languages and text processing utilities (e.g., the widely-used `grep`). The classic algorithm for solving these problems constructs and simulates an NFA corresponding to the regex, which gives an $O(MN)$ running time, where $M$ is the length of the pattern and $N$ is the length of the text.

Unfortunately, significantly faster solutions are unknown and unlikely [6]. However, much faster algorithms exist for many special cases of the problem: dictionary matching; wildcard matching; subset matching; and the word break problem; see [6] and references therein.

Special cases of EDSI have also been studied. First, let us consider the case when both $T_1$ and $T_2$ are degenerate strings. In this case, the problem is trivial: EDSI has a positive answer if and only if for every $i$, $T_1[i] \cap T_2[i]$ is nonempty. Alzamel et al. [2] studied the case when $T_1$ and $T_2$ are generalized degenerate strings: for any $i \in [1, n_1]$ and $j \in [1, n_2]$ all strings in $T_1[i]$ have the same length $\ell_1[i]$, and all strings in $T_2[j]$ have the same length $\ell_2[j]$. In the case of generalized degenerate strings, they showed that deciding if $L(T_1)$ and $L(T_2)$ have a nonempty intersection can be done in $O(N_1 + N_2)$ time. If $T_2$ is a standard string, i.e., an ED string with $m_2 = n_2 = 1$, then we can resort to the results of Bernardini et al. [10] for ED string matching. In particular: there is no combinatorial algorithm\(^2\) for EDSI working in $O(n_1 N_2^{1.5-\epsilon} + N_1)$ time unless the BMM conjecture is false; and we can solve EDSI in $\tilde{O}(n_1 N_2^{1-1} + N_1)$ time. Moreover, Gawrychowski et al. [21] provided a systematic study of the complexity of degenerate string comparison under different notions of matching: Cartesian tree matching; order-preserving matching; and parameterized matching.

Similar to ED strings (and to generalized degenerate strings) is the representation of pangomes via founder graphs. The idea behind founder graphs is that a multiple alignment of few founder sequences can be used to approximate the input MSA, with the feature that each row of the MSA is a recombination of the founders. Like founder graphs, ED strings support the recombination of different rows of the MSA between consecutive columns. Unlike

\(^2\) The notion of “combinatorial algorithm” is informal but widely used in the literature. Typically, we call an algorithm “combinatorial” if it does not not call an oracle for ring matrix multiplication.
ED strings, for which no efficient index is probable [22] (and indeed their value is to enable fast on-line string matching), some subclasses of founder graphs are indexable, and a recent research line is devoted to constructing and indexing such structures [18, 35, 38, 39].

2 Conditional Lower Bounds

In this section, we show several conditional lower bounds for the EDSI problem. Bounds in the first group (see Section 2.1) are based on the popular Strong Exponential-Time Hypothesis (SETH) [11]; the second group of bounds (see Section 2.2) is based on another popular conjecture, the Boolean Matrix Multiplication (BMM) conjecture [1].

2.1 Lower Bounds Based on SETH

We are going to reduce the Orthogonal Vectors (OV, in short) problem to the EDSI problem. In the OV problem we are given a set \( V = \{v^1, \ldots, v^k\} \) of \( k \) binary vectors, each of length \( d \), and we are asked to decide whether or not there are any two vectors in \( V \) which are orthogonal; i.e., the dot product of the two vectors is zero. The OV conjecture, implied by SETH (see [44]), is the following.

**Conjecture 1 (OV conjecture [44]).** The OV problem for \( k \) binary vectors, each of length \( d = \Theta(\log k) \), cannot be solved in \( O(k^{2-\epsilon}) \) time, for any constant \( \epsilon > 0 \).

We show the following reduction.

**Theorem 2.** Given any set \( V = \{v^1, \ldots, v^k\} \) of \( k \) binary vectors of length \( d \), we can construct in linear time two ED strings \( T_1 \) and \( T_2 \) over a binary alphabet such that:

- \( T_1 \) has length, cardinality, and size \( \Theta(ks) \);
- \( T_2 \) has length \( \Theta(\log k) \), cardinality \( \Theta(k) \) and size \( \Theta(ks) \); and
- \( V \) contains two orthogonal vectors if and only if \( T_1 \) and \( T_2 \) have a nonempty intersection.

**Proof.** Let \( u^i = 1^d - v^i \) for all \( i \in [1, k] \). For a length-\( d \) vector \( v \) and \( j \in \{1, \ldots, d\} \), by \( v_j \) we denote the \( j \)th component of \( v \). We construct \( T_1 \) and \( T_2 \) as follows (see Example 3):

\[
T_1 = \prod_{i=1}^{k} \prod_{j=1}^{d} \{0, u^i_j\}, \quad T_2 = \prod_{i=0}^{[\log_2 k]} \{0^{d2^i}, \varepsilon\} \cdot V \cdot \prod_{i=0}^{[\log_2 k]} \{0^{d2^i}, \varepsilon\}.
\]

We now show that \( \mathcal{L}(T_1) \) and \( \mathcal{L}(T_2) \) have a nonempty intersection if and only if there exists a pair of orthogonal vectors in \( V \).

1. Suppose \( v^a \) and \( v^b \) are orthogonal. Then for all \( j \in [1, d] \), \( v^a_j \in \{0, u^j_a\} \) and hence \( v^b \in \prod_j \{0, u^j_a\} \). It follows that

\[
0^{(a-1)d}v^b0^{(k-a)d}d \in \prod_j \{0, u^d_a\} 0^{(a-1)d}d \subseteq \mathcal{L}(T_1).
\]

By decomposing \( a-1 = \sum_{i \in S_{a-1}} 2^i \) and \( k-a = \sum_{i \in S_{k-a}} 2^i \), where for any integer \( p \), the set \( S_p \) contains the positions with a 1 in the binary representation of \( p \), we find that

\[
0^{(a-1)d}v^b0^{(k-a)d}d \in \prod_{i \in S_{a-1}} 0^{d2^i} \cdot V \cdot \prod_{i \in S_{k-a}} 0^{d2^i} \subseteq \mathcal{L}(T_2).
\]

We conclude that \( 0^{(a-1)d}v^b0^{(k-a)d}d \in \mathcal{L}(T_1) \cap \mathcal{L}(T_2) \).

---

3 By the \( \prod \) notation we denote a sequence of concatenations of segments in an ED string.
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- Conversely, suppose that \( \mathcal{L}(T_1) \) and \( \mathcal{L}(T_2) \) have a nonempty intersection and consider a string \( S \in \mathcal{L}(T_1) \cap \mathcal{L}(T_2) \). Let \( v^b \) be the vector from \( V \) which is chosen in \( T_2 \) when constructing \( S \). The strings in the sets of \( T_2 \) all have length divisible by \( d \). Thus \( v^b \) starts at an index \((a-1)d + 1\) of string \( S \) for some integer \( a \). Since \( S \in \mathcal{L}(T_1) \), we have \( v^b \in \prod_{j=1}^{d} \{0, u_j^2\} \). This implies that \( v^a \) and \( v^b \) are orthogonal.

Therefore, solving the orthogonal vectors problem for \( V \) is equivalent to checking whether \( \mathcal{L}(T_1) \) and \( \mathcal{L}(T_2) \) have a nonempty intersection.

**Example 3.** Let \( k = 3, d = 2 \) and \( V = \{v^1 = (1, 0), v^2 = (0, 1), v^3 = (1, 1)\} \).

We have that \( T_1 = \{0\} \cdot \left\{ \begin{array}{c} 0 \\ 1 \end{array} \right\} \cdot \{0\} \cdot \{0\} \cdot \{0\} \) 
and \( T_2 = \left\{ \begin{array}{c} 00 \\ \varepsilon \end{array} \right\} \cdot \left\{ \begin{array}{c} 0000 \\ \varepsilon \end{array} \right\} \cdot \left\{ \begin{array}{c} 10 \\ 01 \end{array} \right\} \cdot \left\{ \begin{array}{c} 00 \\ \varepsilon \end{array} \right\} \cdot \left\{ \begin{array}{c} 0000 \\ \varepsilon \end{array} \right\} \).

One can observe that each string from \( \mathcal{L}(T_1) \cap \mathcal{L}(T_2) \) corresponds to a pair of orthogonal vectors from \( V \). For example, the string \( 010000 \) is in \( \mathcal{L}(T_2) \) because \( v^2 = (0, 1) \in V \). Since the vector \( v^1 = (1, 0) \in V \) is orthogonal to \( v^2 \), one also has \( 010000 \in \mathcal{L}(T_1) \). This is because the two first segments of \( T_1 \) are constructed to encode any vector which is orthogonal to \( v^1 \).

Note that when \( d = \Theta(\log k) \), the length \( n_1 \), the cardinality \( m_1 \) and the size \( N_1 \) of \( T_1 \) are \( O(k \log k) \), whereas \( T_2 \) has length \( n_2 = O(\log k) \), cardinality \( m_2 = O(k) \) and size \( N_2 = O(k \log k) \). Moreover, both ED strings are over a binary alphabet \( \Sigma = \{0, 1\} \). This implies various hardness results for EDSI. For example, we can see that, for any constant \( \epsilon > 0 \), and an alphabet \( \Sigma \) of size at least 2 the problem cannot be solved in \( O((N_1 + N_2 + n_1 + n_2)^{2-\epsilon} \cdot \text{poly}(n_2)) \) time, conditional on the OV conjecture. By using the fact that \( n_1 \leq m_1 \leq N_1 \) and \( n_2 \leq m_2 \leq N_2 \), we obtain the following bounds.

**Corollary 4.** For any constant \( \epsilon > 0 \), there exists no
- \( O((N_1 N_2)^{1-\epsilon})\)-time
- \( O((N_1 m_2 + N_2 m_1)^{1-\epsilon})\)-time
- \( O((N_1 n_2 + N_2 n_1)^{1-\epsilon})\)-time

algorithm for the EDSI problem, unless the OV conjecture is false.

### 2.2 Combinatorial Lower Bounds Based on BMM Conjecture

In the **Triangle Detection (TD)** (in short) problem we are given three \( D \times D \) Boolean matrices \( A, B, C \) and are to check if there are three indices \( i, j, k \in [0, D) \) such that \( A[i,j] = B[j,k] = C[k,i] = 1 \). It is known that Boolean Matrix Multiplication (BMM) and TD either both have truly subcubic combinatorial algorithms or none of them do [45]. The BMM conjecture is stated as follows.

**Conjecture 5** ([BMM conjecture [1]]). Given two \( D \times D \) Boolean matrices, there is no combinatorial algorithm for BMM working in \( O(D^{3-\epsilon}) \) time, for any constant \( \epsilon > 0 \).

Our construction is based on the construction of Bernardini et al. from [10] for ED string matching.

**Theorem 6.** If EDSI over a binary alphabet can be solved in \( O((N_1 + N_2)^{1-2-\epsilon} f(n_1, n_2)) \) time, for any constant \( \epsilon > 0 \) and any function \( f \), then there exists a truly subcubic combinatorial algorithm for TD.
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**Proof.** Let $D$ be a positive integer and let $A$, $B$, and $C$ be three $D \times D$ Boolean matrices. Further let $s \leq D$ be a positive integer to be set later. In the rest of the proof, we can assume that $s$ divides $D$, up to adding $\alpha$ rows and columns containing only 0’s to all three matrices, where $\alpha$ is the smallest non-negative representative of the equivalence class $-D \mod s$.

Let us first construct an ED string $T_1 = X_1X_2X_3$ over a large alphabet with $n_1 = 3$, where each $X_p$, $p \in [1, n_1]$, contains a string for each occurrence of value 1 in $A$, $B$ and $C$, respectively. Below $i$ iterates over $[0, D)$, $j$ and $k$ over $[0, \frac{D}{s})$, and $x$ and $y$ iterate over $[0, s)$. Moreover, $x, y \in [0, s)$, $v_i$ for $i \in [0, D)$, and $a$, $s$ are all letters.

- If $A[i, x \cdot \frac{D}{s} + j] = 1$, then $X_1$ contains the string $v_i xa^j$;
- If $B[x \cdot \frac{D}{s} + j, y \cdot \frac{D}{s} + k] = 1$, then $X_2$ contains the string $a^{\frac{D}{s} - j}x\$y$a^{\frac{D}{s} - k}$;
- If $C[y \cdot \frac{D}{s} + k, i] = 1$, then $X_3$ contains the string $a^i y v_i$;

The length of each string in each $X_p$ is $O(D/s)$ and the total number $m_1$ of strings is up to $3D^2$. Overall, $N_1 = O(D^3/s)$.

We construct an ED string $T_2$ with $n_2 = 1$ containing the following strings:

$$P(i, x, y) = v_i x_a^2 y_a^2 y_v_i \quad \text{for every } x, y \in [0, s) \text{ and } i \in [0, D).$$

Each string has length $O(D/s)$ and there are $m_2 = Ds^2$ strings, so $N_2 = O(D^2 s)$.

We use the following fact.

**Fact 7** ([10]). $P(i, x, y) \in \mathcal{L}(T_1)$ if and only if the following holds for some $j, k \in [0, D/s)$:

$$A[i, x \cdot \frac{D}{s} + j] = B[x \cdot \frac{D}{s} + j, y \cdot \frac{D}{s} + k] = C[y \cdot \frac{D}{s} + k, i] = 1.$$

We choose $s = \lceil \sqrt{D} \rceil$; then $N_1, N_2 = O(D^2)$ and $n_1, n_2 = O(1)$. Then indeed an $O((N_1 + N_2)^{1.5 - \epsilon} \phi(n_1, n_2))$-time algorithm for EDSI would yield an $O(D^{3 - 2\epsilon})$-time algorithm for the TD problem.

Note also that even though the size of the alphabet used above is $\Theta(s + D) = \Theta(D)$, we can encode all letters by equal-length binary strings blowing $N_1$ and $N_2$ up only by a factor of $\Theta(\log D)$ and, hence, obtain the same lower bound for a binary alphabet.

Both $m_1$ and $m_2$ in the reduction are $O(D^2)$, thus an $O((m_1 + m_2)^{1.5 - \epsilon} \phi(n_1, n_2))$-time algorithm would yield an $O(D^{3 - 2\epsilon})$-time algorithm for TD; hence we obtain the following.

**Corollary 8.** If EDSI over a binary alphabet can be solved in $O((N_1^{1.2} + N_2^{1.2} + m_1^{1.5} + m_2^{1.5})^{1 + \epsilon} \phi(n_1, n_2))$ time, for any constant $\epsilon > 0$ and any function $\phi$, then there exists a truly subcubic combinatorial algorithm for TD.

## 3 EDSI: The Unary Case

An ED string is called *unary* if it is over an alphabet of size 1. In this special case, if both $T_1$ and $T_2$ are over the same alphabet $\Sigma = \{a\}$, EDSI boils down to checking whether there exists any $b \geq 0$ such that $a^b$ belongs to both $\mathcal{L}(T_1)$ and $\mathcal{L}(T_2)$.

Let $T$ be a unary ED string of length $n$ over alphabet $\Sigma = \{a\}$. We define the *compact representation* $R(T)$ of $T$ as the following sequence of sets of integers:

$$\forall i \in [1, n] \quad R(T)[i] = \{b_{i,1}, b_{i,2}, \ldots, b_{i,m_i}\} \iff T[i] = \{a^{b_{i,1}}, a^{b_{i,2}}, \ldots, a^{b_{i,m_i}}\},$$

where $b_{i,j} \geq 0$ for all $i \in [1, n]$ and $j \in [1, m_i]$, the cardinality of $T$ is $m = \sum_{i=1}^{n} m_i$, and its size is $N = N_\epsilon + \sum_{i=1}^{n} \sum_{j=1}^{m_i} b_{i,j}$, where $N_\epsilon$ is the total number of empty strings in $T$. 
Comparing Elastic-Degenerate Strings: Algorithms, Lower Bounds, and Applications

Theorem 9. If $T_1$ and $T_2$ are unary ED strings and each is given in a compact representation, the problem of deciding whether $\mathcal{L}(T_1) \cap \mathcal{L}(T_2)$ is nonempty is NP-complete.

Proof. The problem is clearly in NP, as it is enough to guess a single element for each set in both $T_1$ and $T_2$, and then simply check if the sums match in linear time. We show the NP-hardness through a reduction from the Subset Sum problem, which takes $n$ integers $b_1, b_2, \ldots, b_n$ and an integer $c$, and asks whether there exist $x_i \in \{0, 1\}$, for all $i \in [1, n]$, such that $\sum_{i=1}^{n} x_i b_i = c$. Subset Sum is NP-complete [30] also for non-negative integers. For any instance of Subset Sum, we set $R(T_1)[i] = \{b_i, 0\}$ for all $i \in [1, n]$, $n_2 = 1$ and $R(T_2)[1] = \{c\}$. Then the answer to the Subset Sum instance is YES if and only if $\mathcal{L}(T_1) \cap \mathcal{L}(T_2)$ is nonempty.

In what follows, we provide an algorithm which runs in polynomial time in the size of the two unary ED strings when the latter are given uncompacted.

The set $\mathcal{L}(T)$ can be represented as a set $L(T) \subset \mathbb{N}$ such that $\mathcal{L}(T) = \{a^\ell : \ell \in L(T)\}$. The set $L(T)$ will be stored as a list (without repetitions). We will show how to efficiently compute $L(T_1)$ and $L(T_2)$. Then one can compute $L(T_1) \cap L(T_2)$ in $O(N_1 + N_2)^2$ time, which allows, in particular, to check if $L(T_1) \cap L(T_2) = \emptyset$ (which is equivalent to $\mathcal{L}(T_1) \cap \mathcal{L}(T_2) = \emptyset$).

We show the computation for $L(T_1)$. The workhorse is an algorithm from the following Lemma 10 that allows to compute the set $L(X_1X_2)$ of concatenation of two ED strings based on their sets $L(X_1), L(X_2)$.

Lemma 10. Let $X_1$ and $X_2$ be ED strings. Given $L(X_1)$ and $L(X_2)$ such that $t_1 = \max L(X_1)$ and $t_2 = \max L(X_2)$, we can compute $L(X_1X_2)$ in $O((t_1 + t_2)^2 \log(t_1 + t_2))$ time.

Proof. For two sets $A, B \subset \mathbb{N}$, by $A + B$ we denote the set $\{a + b : a \in A, b \in B\}$. We then have $L(X_1X_2) = L(X_1) + L(X_2)$. Fast Fourier Transform (FFT) [14] can be used directly to compute $L(X_1) + L(X_2)$ in $O((t_1 + t_2)^2 \log(t_1 + t_2))$ time.

Lemma 11. $L(T_1)$ can be computed in $O(N_1 \log N_1 \log n_1)$ time.

Proof. We apply the recursive algorithm described in Algorithm 1 to $T_1$.

Algorithm 1 Compute-$L(T[1] \cdots T[k])$.

if $k = 1$ then
  Compute $L(T[1])$ naively
  $i$ ← $\lfloor k/2 \rfloor$
  $L_1$ ← Compute-$L(T[1] \cdots T[i])$
  $L_2$ ← Compute-$L(T[i+1] \cdots T[k])$
  return $L_1 + L_2$

Let $N_{i,i} = \sum_{x \in \mathcal{L}(T[i])} x$ and $t_{i,i} = \max L(T[i])$ for $i \in [1, n_1]$. Obviously, $t_{i,i} \leq N_{i,i}$.

We analyze the complexity of the recursion by levels. For the bottom level, $L(T[1][i])$ can be computed in $O(N_{i,i})$ time for each $i \in [1, n_1]$, which sums up to $O(N_1)$. For the remaining levels, we notice that $\max L(T_1[i] \cdots T_1[j]) = t_{1,i} + \cdots + t_{1,j}$. On each level, the fragments of $T_1$ that are considered are disjoint. Thus, the complexity on each level via Lemma 10 is $O((\sum_{i=1}^{n_1} t_{1,i}) \log(\sum_{i=1}^{n_1} t_{1,i})) = O(N_1 \log N_1)$. The number of levels of recursion is $O(\log n_1)$; the complexity follows.

\[
\]
Theorem 12. If $T_1$ and $T_2$ are unary ED strings, then $L(T_1) \cap L(T_2)$ can be computed in $O(N_1 \log N_1 \log n_1 + N_2 \log N_2 \log n_2)$ time.

Proof. We use Lemma 11 to compute $L(T_1)$ and $L(T_2)$ in the required complexity. Then $L(T_1) \cap L(T_2)$ can be computed via bucket sort.

4 EDSI: General Case

Assuming that the two ED strings, $T_1$ and $T_2$, of total size $N_1 + N_2$ are over an integer alphabet $[1, (N_1 + N_2)^{O(1)}]$, we can sort the suffixes of all strings in $T_1[i]$, for all $i \in [1, n_1]$, and the suffixes of all strings in $T_2[j]$, for all $j \in [1, n_2]$, in $O(N_1 + N_2)$ time [19].

By LCP($X, Y$) let us denote the length of the longest common prefix of two strings $X$ and $Y$. Given a string $S$ over an integer alphabet, we can construct a data structure over $S$ in $O(|S|)$ time, so that when $i, j \in [1, |S|]$ are given to us on-line, we can determine LCP($S[i \ldots |S|], S[j \ldots |S|]$) in $O(1)$ time [15].

4.1Compacted NFA Intersection

In this section we show an algorithm for computing a representation of the intersection of the languages of two ED strings using techniques from formal languages and automata theory.

Definition 13 (NFA). A nondeterministic finite automaton (NFA) is a 5-tuple $(Q, \Sigma, \delta, q_0, F)$, where $Q$ is a finite set of states; $\Sigma$ is an alphabet; $\delta : Q \times (\Sigma \cup \{\varepsilon\}) \rightarrow P(Q)$ is a transition function, where $P(Q)$ is the power set of $Q$; $q_0 \in Q$ is the starting state; and $F \subseteq Q$ is the set of accepting states.

Using the folklore product automaton construction, one can check whether two NFA have a nonempty intersection in $O(N_1 \cdot N_2)$ time, where $N_1$ and $N_2$ are the sizes of the two NFA [33]. We use a different, compacted representation of automata, which in some special cases allows a more efficient algorithm for computing and representing the intersection.

Definition 14 (Compacted NFA). An extended transition is a transition function of the form $\delta^{ext} : Q \times \Sigma^* \rightarrow P(Q)$, where $Q$ is a finite set of states, $\Sigma^*$ is the set of strings over alphabet $\Sigma$, and $P(Q)$ is the power set of $Q$. A compacted NFA is an NFA in which we allow extended transitions. Such an NFA can also be represented by a standard (uncompacted) NFA, where each extended transition is subdivided into standard one-letter transitions (and $\varepsilon$-transitions), $\delta : Q \times (\Sigma \cup \{\varepsilon\}) \rightarrow P(Q)$. The states of the compacted NFA are called explicit, while the states obtained due to these subdivisions are called implicit.

Given a compacted NFA $A$ with $V$ explicit states and $E$ extended transitions, we denote by $V^e$ and $E^e$ the number of states and transitions, respectively, of its uncompacted version $A^u$. Henceforth we assume that in the given NFA every state is reachable, and hence we have $V^e = O(E^e)$ and $V = O(E)$.

Lemma 15. Given two compacted NFA $A_1$ and $A_2$, with $V_1$ and $V_2$ explicit states and $E_1$ and $E_2$ extended transitions, respectively, a compacted NFA representing the intersection of $A_1$ and $A_2$ with $O(V_1^e V_2 + V_1 V_2^e)$ explicit states and $O(E_1^e E_2 + E_1 E_2^e)$ extended transitions can be computed in $O(E_1^e + E_2^e)$ time.

Proof. We start by constructing an LCP data structure over the concatenation of all the labels of extended transitions of both NFA of total size $O(E_1^e + E_2^e)$. It requires $O(E_1^e + E_2^e)$-time preprocessing and allows answering LCP queries on any two substrings of such labels in $O(1)$ time.
We construct \( B \) a compacted NFA representing the intersection of \( A_1 \) and \( A_2 \).

Every state of \( B \) is composed of a pair: an explicit state of one automaton and any explicit or implicit state of the other automaton (or equivalently a state of the uncompiled version of the automaton). Thus the total number of explicit states of \( B \) is \( O(V_1^n V_2 + V_1 V_2^n) \).

We need to compute the extended transitions of \( B \). For a state \((u, v)\) we check every string pair \((P, Q)\), where \( P \) iterates over all extended transitions going out of \( u \) and \( Q \) iterates over all extended transitions going out of \( v \) (a transition going out of an implicit state is represented by a suffix of the transition it belongs to). For every pair \((P, Q)\) we ask an \( \text{LCP}(P, Q) \) query. If \( \text{LCP}(P, Q) \) is equal to one of \(|P|, |Q|\) (possibly both), we create an extended transition between \((u, v)\) and the pair of states reachable through those transitions (if one of the transitions is strictly longer, we prune it to the right length, ending it at an implicit state of its input NFA). Otherwise such a transition does not lead to any explicit state of \( B \) and thus cannot be used to reach the accepting state; hence we ignore it.

Finally, the starting (resp. accepting) state of \( B \) corresponds to a pair of starting (resp. accepting) states of \( A_1 \) and \( A_2 \).

Since any pair representing an explicit state of \( B \) contains an explicit state of \( A_1 \) or \( A_2 \), the number of such transition pair checks (and hence also the number of the extended transitions of \( B \)) is \( O(E_1^n E_2 + E_1 E_2^n) \). Since each such check takes \( O(1) \) time, the construction complexity follows. Note that NFA \( B \) may contain unreachable states; such states can be removed afterwards in linear time. The algorithms’ correctness follows from the observation that \( B^n \) is in fact the standard intersection automaton of \( A_1^n \) and \( A_2^n \) with some states, that do not belong to any path between the starting and the accepting states, removed.

We next define the path-automaton of an ED string (inspect Figure 3 for an example).

**Definition 16 (Path-automaton).** Let \( T \) be an ED string of length \( n \), cardinality \( m \), and size \( N \). The path-automaton of \( T \) is the compacted NFA consisting of:

- \( V = n + 1 \) explicit states, numbered from 1 through \( n + 1 \). State 1 is the starting state and state \( n + 1 \) is the accepting state. State \( i \in [2, n] \) is the state in-between \( T[i−1] \) and \( T[i] \).
- \( m_i \) extended transitions from state \( i \) to state \( i + 1 \) labeled with the strings in \( T[i] \), for all \( i \in [1, n] \), where \( E = m = \sum_i m_i \).

The path-automaton of \( T \) accepts exactly \( L(T) \). The uncompacted version of this path-automaton has \( V^n = O(N) \) states and \( E^n = N \) transitions.

Lemma 15 thus implies the following result.

**Corollary 17.** The compacted NFA representing the intersection of two path-automata with \( O(N_1 m_2 + N_2 m_1) \) explicit states and \( O(N_1 m_2 + N_2 m_1) \) extended transitions can be constructed in \( O(N_1 m_2 + N_2 m_1) \) time.
Theorem 18. EDSI can be solved in \(O(N_1m_2 + N_2m_1)\) time. If the answer is YES, a witness can be reported within the same time complexity.

Proof. The path-automaton of an ED string of size \(N\) can be constructed in \(O(N)\) time. Given two ED strings, we can construct their path-automata in linear time and apply Corollary 17. By finding any path from the starting to the accepting state in linear time (if it exists), we obtain the result.

Notice that the path-automata representing ED strings, as well as their intersection, are always acyclic, but may contain \(\varepsilon\)-transitions. In the following we are only interested in the graph underlying the path-automaton, that is the directed acyclic graph (DAG), where every node represents an explicit state and every labeled directed edge represents an extended transition of the path-automaton (inspect also Figure 1).

4.2 An \(\tilde{O}(N_1^{\omega - 1}n_2 + N_2^{\omega - 1}n_1)\)-time Algorithm for EDSI

In this section, we start by showing a construction of the intersection graph computed by means of Lemma 15 in the case when the input is a pair of path-automata that allows an easier and more efficient implementation. The construction is then adapted to obtain an \(\tilde{O}(N_1^{\omega - 1}n_2 + N_2^{\omega - 1}n_1)\)-time algorithm for solving the EDSI problem.

For \(x \in \{1, 2\}\) by \(A_x\) we denote the compacted NFA (henceforth, graph \(A_x\)) representing the ED string \(T_x\). By \(I_x[i]\) we denote the set of implicit states (henceforth, implicit nodes) appearing on the extended transitions (henceforth, edges) between explicit states (henceforth, explicit nodes) \(i\) and \(i + 1\). For convenience, the implicit nodes in the sets \(I_x[1], \ldots, I_x[n_x]\) can be numbered consecutively starting from \(n_x + 2\).

Let \(U_{i,j} = \{(i, k) : k \in \{j\} \cup I_x[j]\}\) and \(U'_{i,j} = \{(k, j) : k \in \{i\} \cup I_x[i]\}\), for all \(i \in [1, n_1 + 1]\) and \(j \in [1, n_2 + 1]\). As in the construction of Lemma 15, the union of all \(U_{i,j}\) and \(U'_{i,j}\) is the set of explicit nodes of the intersection graph that we construct; this can be represented graphically by a grid, where the horizontal and vertical lines correspond to \(U_{i,j}\) and \(U'_{i,j}\), respectively (inspect Figure 4a). In particular, we would like to compute the edges between these explicit nodes (inspect Figure 4b) in \(O(N_1m_2 + N_2m_1)\) time.

Consider an explicit node of the intersection graph; this node is represented by a pair of nodes: one from \(A_1\) and one from \(A_2\). We need to consider two cases: explicit vs explicit node; or explicit vs implicit node. Without loss of generality, we consider the first node to be explicit. Let us denote this pair by \((i, k) \in U_{i,j}\), where \(i\) is an explicit node of \(A_1\) and \(k\) is a node of \(A_2\). Let us further denote by \(\ell_1\) the label of one of the edges going from node \(i\) to node \(i + 1\). For \(k\), we have two cases. If \(k\) is explicit (i.e., \(k = j\)) then we denote by \(\ell_2\) the label of one of the edges going from \(k\) to \(k + 1\). Otherwise (\(k\) is implicit), we denote by \(\ell_2\) the path label (concatenation of labels) from node \(k\) to node \(j + 1\).

As noted in the proof of Lemma 15, an edge is constructed only if \(\operatorname{LCP}(\ell_1, \ell_2) = \min(|\ell_1|, |\ell_2|)\). If \(\operatorname{LCP}(\ell_1, \ell_2) = |\ell_2| < |\ell_1|\) (a prefix of a string in \(T_1[i]\) is equal to the suffix of a string in \(T_2[j]\) starting at the position corresponding to node \(k \in \{j\} \cup I_x[j]\)), the edge ends in a node from \(U'_{i,j+1}\) (Figure 4c). If \(\operatorname{LCP}(\ell_1, \ell_2) = |\ell_1| < |\ell_2|\) (a whole string from \(T_1[i]\) occurs in a string from \(T_2[j]\) starting at the position corresponding to node \(k \in \{j\} \cup I_x[j]\)), the edge ends in a node from \(U_{i+1,j}\) (Figure 4d). Otherwise \(\operatorname{LCP}(\ell_1, \ell_2) = |\ell_1| = |\ell_2|\); the two strings are equal) the edge ends in \((i + 1, j + 1)\). Symmetrically (i.e., the second node is explicit), the edge going out of a node from \(U'_{i,j}\) ends at a node from the same set \(U'_{i,j+1} \cup U_{i+1,j} \cup \{(i + 1, j + 1)\}\) (inspect Figure 4b).
Figure 4 An overview of the edges computed by the algorithm.

We next show how to construct the intersection graph by computing all such edges going out of $U_{i,j}$ or $U'_{i,j}$ in a single batch using suffix trees (inspect Figure 5 in Appendix A for an example). This construction allows an easier and more efficient implementation in comparison to the LCP data structure used in the general NFA intersection construction. Let us recall that $||T||$ denotes the size of the ED string $T$.

**Lemma 19.** For any $i \in [1, n_1 + 1]$ and $j \in [1, n_2 + 1]$, we can construct all $K_{i,j}$ edges going out of nodes in $U_{i,j}$ in $O(N_{1,i} + N_{2,j} + K_{i,j})$ time, where $N_{1,i} = ||T_1[i]||$ and $N_{2,j} = ||T_2[j]||$, using the generalized suffix tree of the strings in $T_2[j]$.

**Proof.** We first construct the generalized suffix tree of the strings in $T_2[j]$ in $O(N_{2,j})$ time [19]. We also mark each node corresponding to a suffix of a string in $T_2[j]$ with a $T_2$-label. Each such node is also decorated with one or multiple starting positions, respectively, from one or multiple elements of $T_2[j]$ sharing the same suffix. For each branching node of the suffix tree, we construct a hash table, to ensure that any outgoing edge can be retrieved in constant time based on the first letter (the key) of its label. This can be done in $O(N_{2,j})$ time with perfect hashing [20]. We next spell each string from $T_1[i]$ from the root of the suffix tree
making implicit nodes explicit or adding new ones if necessary to create the compacted trie of all those strings; and, finally, we mark the reached nodes of the suffix tree with a \( T_1 \)-label. Spelling all strings from \( T_1[i] \) takes \( O(N_{1,i}) \) time.

Every pair of different labels marking two nodes in an ancestor-descendant relationship corresponds to exactly one outgoing edge of the nodes in \( U_{i,j} \): (i) if a node marked with a \( T_2 \)-label is an ancestor of a node marked with a \( T_1 \)-label, then the suffix of a string from \( T_2[j] \) matches a prefix of a string from \( T_1[i] \) forming an edge ending in \( U'_{i,j+1} \); (ii) if a node marked with a \( T_1 \)-label is an ancestor of a node marked with a \( T_2 \)-label, then a string from \( T_1[i] \) occurs in a string from \( T_2[j] \) extending its prefix and forming an edge ending in \( U_{i+1,j} \); (iii) if a node is marked with a \( T_1 \)-label and with a \( T_2 \)-label, then the suffix of a string from \( T_2[j] \) matches a string from \( T_1[i] \) forming an edge ending in \((i + 1, j + 1)\). After constructing the generalized suffix tree of \( T_2[j] \) and spelling the strings from \( T_1[i] \), it suffices to make a DFS traversal on the annotated tree to output all \( K_{i,j} \) such pairs of nodes.

\[ \textbf{Theorem 20.} \text{We can construct the intersection graph of } T_1 \text{ and } T_2 \text{ in } O(N_1n_2 + N_2m_1) \text{ time using the suffix tree data structure and tree search traversals.} \]

\textbf{Proof.} We apply Lemma 19 for \( U_{i,j} \) and \( U'_{i,j} \), for all \( i \in [1, n_1 + 1] \) and \( j \in [1, n_2 + 1] \). We have that the total number of nodes is \( \sum_{i,j} O(N_{1,i} + N_{2,j}) = O(N_1n_2 + N_2m_1) \), and then the sum of all output edges is bounded by \( O(N_1m_2 + N_2m_1) \) by Corollary 17.

Note that if we are interested only in checking whether the intersection is nonempty, and not in the computation of its graph representation, it suffices to check which of the nodes are reachable from the starting node, which may be more efficient as there are \( O(N_1n_2 + N_2m_1) \) explicit nodes in this graph.

Let \( X \) be the set of nodes of \( U_{i,j} \) that are reachable from the starting node. From this set of nodes we need to compute two types of edges (inspect Figure 4b). The first type of edges, namely, the ones from \( X \) to \( U'_{i,j+1} \cup \{(i + 1, j + 1)\} \) (green edges in Figure 4b) are computed by means of Lemma 21, which is similar to Lemma 19. For the second type of edges, namely, the ones from \( X \) to \( U_{i+1,j} \cup \{(i + 1, j + 1)\} \) (blue edges in Figure 4b), we use a reduction to the so-called active prefixes extension problem \[10\] (Lemma 23).

\[ \textbf{Lemma 21.} \text{For any given } X \subseteq U_{i,j}, \text{ we can compute the subset of } U'_{i,j+1} \cup \{(i + 1, j + 1)\} \text{ containing all and only the nodes that are reachable from the nodes of } X \text{ in } O(N_{1,i} + N_{2,j}) \text{ time.} \]

\textbf{Proof.} In Lemma 19, the edges from nodes of \( U_{i,j} \) to nodes of \( U'_{i,j+1} \) come from a pair of nodes in the generalized suffix tree of \( T_2[j] \): one marked with a \( T_1 \)-label and its ancestor marked with a \( T_2 \)-label. Notice that the \( T_2 \)-labels are in correspondence with the elements of \( U_{i,j} \) (the labels on a proper suffix of a string in \( T_2 \) are in a one-to-one correspondence with \( U_{i,j} \setminus \{(i, j)\} \)), and \( (i, j) \) corresponds to whole strings in \( T_2[j] \), and hence we can trivially remove the \( T_2 \)-labels that do not correspond to the elements of \( X \). Furthermore, we are not interested in the set of starting positions decorating a node with a \( T_2 \)-label; we are interested only in whether a node is \( T_2 \)-labeled or not (i.e., we do not care from which node of \( X \) the edge originates). Since the nodes marked with a \( T_1 \)-label have in total \( N_{1,i} \) ancestors (including duplicates), we can compute the result of this case in \( O(N_{1,i} + N_{2,j}) \) time in total.

Finally, the node \((i + 1, j + 1)\) is reachable when a single node is marked with both a \( T_1 \)-label and a \( T_2 \)-label. This can be checked within the same time complexity.

\footnote{Our implementation of this algorithm can be found at https://github.com/urbanslug/junctions.}
The remaining edges (blue edges in Figure 4b) are dealt with via a reduction to the following problem:

<table>
<thead>
<tr>
<th>Active Prefixes Extension (APE)</th>
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<tbody>
<tr>
<td><strong>Input:</strong> A string $P$ of length $m$, a bit vector $W$ of size $m$, and a set $S$ of strings of total length $N$.</td>
</tr>
<tr>
<td><strong>Output:</strong> A bit vector $V$ of size $m$ with $V[p] = 1$ if and only if there exists $P' \in S$ and $p' \in [1, m]$, such that $P[1..p' - 1] \cdot P' = P[1..p - 1]$ and $W[p'] = 1$.</td>
</tr>
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Bernardini et al. have shown the following result in [10].

- **Lemma 22 ([10]).** The APE problem can be solved in $\tilde{O}(m^{\omega} - 1) + O(N)$ time, where $\omega$ is the matrix multiplication exponent.

- **Lemma 23.** For any given $X \subseteq U_{i,j}$, we can compute the subset of $U_{i+1,j} \cup \{i+1, j+1\}$ containing all and only the nodes that are reachable from the nodes of $X$ in $\tilde{O}(N_{i,i} + N_{2,j})$ time.

**Proof.** The problems of computing the subset of $U_{i+1,j}$ reachable from $X$ and the APE problem can be reduced to one another in linear time.

For the forward reduction, let us set $S = T_1[i]$ and $P = \prod_{S \in T_2[j]} S$, where $S$ is a letter outside of the alphabet of $T_1$. This means that we order the strings in $T_2[j]$, in an arbitrary but fixed way. For a single string $S$ (where $S \in T_2[j]$), the positions from $S[1..|S| - 1]$ correspond to the implicit nodes (along the path spelling $S$) of $I_2[j]$, while the position with $S$ corresponds to the explicit node $j$ of $A_2$ and the one with $S[|S|]$ to the explicit node $j + 1$ of $A_2$. Through this correspondence, we can construct two bit vectors $W$ and $V$, each of them of size $|P|$, and whose positions are in correspondence with $\{j\} \cup I_2[j] \cup \{j + 1\}$ (note that this correspondence is not a bijection, as the explicit nodes $j$ and $j + 1$ have several preimages when $|T_2[j]| \geq 2$). As $U_{i,j} \cup \{(i, j + 1)\}$ and $U_{i+1,j} \cup \{(i+1, j+1)\}$ are copies of $\{j\} \cup I_2[j] \cup \{j + 1\}$, we use the same correspondence to match positions between $W$ and $U_{i,j} \cup \{(i, j + 1)\}$ and between $V$ and $U_{i+1,j} \cup \{(i+1, j+1)\}$. Finally, we set $W[k] = 1$ if and only if the corresponding node of $U_{i,j}$ belongs to $X$ (for $k$ corresponding to $(i,j+1)$, we set $W[k] = 0$ as such a node cannot belong to $X$). After solving APE, we have $V[k] = 1$ for some $k$ corresponding to a node of $U_{i+1,j} \cup \{(i+1, j+1)\}$ if and only if this node is reachable from $X$.

In more detail, observe that since $S$ does not belong to the alphabet of $T_1$, a string $S$ from $T_1[i]$ has to match a fragment of a string from $T_2[j]$ to set $V[k]$ to 1. This happens only if additionally $W[k - |S|] = 1$; both things happen at the same time exactly when: (i) there exists a node $(i, \ell) \in X$; (ii) there exists an edge from $(i, \ell)$ to $(i+1, \ell')$; and (iii) the positions $k - |S|$ and $k$ in $P$ correspond to $\ell, \ell'$, respectively.

In the above reduction we have $|P| = \sum_{S \in T_2[j]} |S| + 1 = O(N_{2,j})$, and $|S| = N_{1,j}$, hence the lemma statement follows by Lemma 22.

For the reverse reduction, given an instance of APE, we encode it by setting $T_1[i] = S, T_2[j] = \{P\}$, $(N_{1,i} = |S|, N_{2,j} = |P|)$ and $X$ containing the nodes corresponding to positions $k$ where $W[k] = 1$ (the last element of such $X$ is potentially $(i+1,j)$, but we do not care about this corner case of extending the prefix which is already the full string $P$).

This reduction shows that a more efficient solution to the problem of finding the endpoints of edges originating in $X$ would result in a more efficient solution to the APE problem. ▶

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5 Here, note that if the node is $(i+1,j)$ or $(i+1,j+1)$, then a corresponding $k$ is not unique, but at least one of them satisfy $V[k] = 1$. 

Theorem 24. We can solve EDSI in $\tilde{O}(N_1^{\omega-1}n_2 + N_2^{\omega-1}n_1)$ time, where $\omega$ is the matrix multiplication exponent. If the answer is YES, we can output a witness within the same time complexity.

Proof. It suffices to set the starting node $(1, 1)$ as reachable, apply Lemmas 21 and 23, and their symmetric versions for $U'_{i,j}$, for each value of $(i, j) \in [1, n_1 + 1] \times [1, n_2 + 1]$ in lexicographical order, with $X$ equal to the set of reachable nodes of $U_{i,j}$ (respectively of $U'_{i,j}$); and, finally, check whether node $(n_1 + 1, n_2 + 1)$ is set as reachable. We bound the total time complexity of the algorithm by:

$$\sum_{i,j} \tilde{O}(N_1^{\omega-1} + N_2^{\omega-1}) = \tilde{O}(n_2 \sum_i N_1^{\omega-1} + n_1 \sum_j N_2^{\omega-1}) \leq \tilde{O}(N_1^{\omega-1}n_2 + N_2^{\omega-1}n_1).$$

If $L(T_1) \cap L(T_2)$ is nonempty, that is, if the node $(n_1 + 1, n_2 + 1)$ is set as reachable from node $(1, 1)$, then we can additionally output a witness of the intersection – a single string from $L(T_1) \cap L(T_2) –$ within the same time complexity. To do that we mimic the algorithm on the graph with reversed edges. This time, however, we do not mark all of the reachable nodes; we rather choose a single one that was also reachable from $(1, 1)$ in the forward direction. This way, the marked nodes form a single path from $(1, 1)$ to $(n_1 + 1, n_2 + 1)$. The witness is obtained by reading the labels on the edges of this path.

5 Acronym Generation

In this section we study a problem on standard strings. Given a sequence $P = P_1, \ldots, P_n$ of $n$ strings we define an acronym of $P$ as a string $A = A_1 \cdots A_n$, where $A_i$ is a (possibly empty) prefix of $P_i$, $i \in [1, n]$. We next formalize the Acronym Generation problem.

**ACRONYM GENERATION (AG)**

**Input:** A set $D$ of $k$ strings of total length $K$ and a sequence $P = P_1, \ldots, P_n$ of $n$ strings of total length $N$.

**Output:** YES if some acronym of $P$ is an element of $D$, NO otherwise.

The AG problem is underlying real-world information systems (e.g., see https://acronymify.com/ or https://acronym-generator.com/) and existing approaches rely on brute-force algorithms or heuristics to address different variants of the problem [41, 40, 32, 34, 28, 43, 29, 31]. These algorithms usually accept a sequence $P$ of $n \leq n_{\text{max}}$ strings, for some small integer $n_{\text{max}}$, which highlights the lack of efficient exact algorithms for generating acronyms. Here we show an exact polynomial-time algorithm to solve AG for any $n$.

We can encode AG by means of EDSI and modify the developed methods. Let $T_1[i]$, $i \in [1, n]$, be the set of all prefixes of $P_i$ and further let $T_2[1] = D$. By using Theorem 18 or Corollary 17 we obtain an $O(\sum_i |P_i|^2k + KN) = O(N^2k + KN)$-time algorithm, while using Theorem 24 we obtain an $\tilde{O}(N^{2\omega-2} + K^{\omega-1}n)$-time algorithm, for solving the AG problem.

Since, however, all elements of set $T_1[i]$ are prefixes of a single string ($P_i$), we can obtain a more efficient graph representation of $T_1$ by joining nodes $i$ and $i + 1$ with a single path labeled with $P_i$, with an additional $\epsilon$ edge between every (implicit) node of the path and node $i + 1$. As the size of the graph for $T_1$ is smaller ($O(N)$ nodes and edges), by using Lemma 15 we obtain an $O(Nk + KN) = O(NK)$-time algorithm for solving the AG problem.

The considered ED strings have additional strong properties however. $T_1[i]$’s are not just sets of prefixes of single strings, but sets of all their prefixes, while the length $n_2$ of $T_2$ is equal to 1. By employing these two properties we obtain the following improved result.
Theorem 25. AG can be solved in $O(nK + N)$ time.

Proof. The algorithm of Theorem 24 is based on finding out which elements of sets $U_{i,j}, U'_{i,j}$ are reachable; however, since $n_2 = 1$, the sets $U'_{i,j}$ are trivialized: by definition, a node from the middle of $T_1[i]$ cannot correspond to the starting or accepting node of the graph of $T_2$ (reading a letter in the first graph means also moving out of the starting node in the second one), hence the only possible reachable node of $U'_{i,j}$ is the explicit node $(i, j)$, which is also an element of $U_{i,j}$. More formally, a reachable node $(k, 1) \in U'_{i,1}$ must be equal to $(i, 1)$ as other such nodes can only be reached using some edge with a nonempty label. By symmetry, nodes from $U'_{1,j}$ other than $(i, 2)$ are not backwards reachable from the accepting node.

In Lemma 23, to compute the reachable nodes of $U_{i+1,j}$ knowing the reachable nodes of $U_{i,j}$, fast matrix multiplication is employed (Lemma 22), but in this special case a simpler method will be more effective. Let $W_k$ be the string read between nodes $k$ and 2 in the path-graph of $T_2$. The crucial observation is: the edges going out of node $(i, k) \in U_{i,1} \cup \{(i, 2)\}$ for $k \neq 1$ end in nodes $(i + 1, k')$ for $k' \in [k, k + l]$, where $l = \text{LCP}(P_i, W_k)$ as the strings from $T_1[i]$ matching the prefix of $W_k$ are exactly all the prefixes of $P_i$ of length at most $l$.

Hence to compute the reachable subset of $U_{i+1,1} \cup \{(i + 1, 2)\}$, we can handle the edges going out of $(i, 1)$ separately in $O(K + |P_i|)$ time by letter comparisons, then compute the LCP($P_i, W_k$) for all the reachable nodes $(i, k)$ either using the LCP data structure, or with the use of the generalized suffix tree of $T_2[i] = D$ in $O(K + |P_i|)$ total time, and finally, using a sweep line approach, compute the union of the obtained intervals in $O(K)$ time. We answer YES if and only if node $(n + 1, 2)$ is reachable.

Over all $i$ values this gives an algorithm running in $\sum_i O(K + |P_i|) = O(nK + N)$ total time. Furthermore one is allowed to choose, for each $i \in [1, n]$, the minimal length $x_i$ of the prefix of $P_i$ (including length $x_i = 0$ if one wants to allow empty prefixes) used in the acronym (some strings should not be completely excluded from the acronym). The only modification to the algorithm in such a generalized case is replacing intervals $[k, k + l]$ by $[k + x_i, k + l]$, which does not influence the claimed complexity.

Corollary 26. If the answer to the instance of the AG problem is YES, we can output all strings in $D$ which are acronyms of $P$ within $O(nK + N)$ time.

Proof. In the algorithm employed by Theorem 25 the reachable nodes of $U_{i,1} \cup \{(i, 2)\}$ are found. When the node $(i + 1, 2)$ is the endpoint of an edge starting in node $(i, k)$ for $k \neq 1$, then the path of the path-graph of $T_2$ containing node $k$ is an acronym of $P$. If node $(i + 1, 2)$ is reached directly from reachable node $(i, 1)$, then the whole prefix of $P_i$ used to do that is in $D$, and hence is a standalone acronym of $P$. If for a path neither of the two cases qualifies, then it cannot be used to reach node $(n + 1, 2)$, and hence is not an acronym of $P$.

If the generalization with minimal lengths of prefixes is applied, then the values of $i$ used here are restricted to $[i', n]$, where $i'$ is the largest value of $i$ with a restriction $x_i > 0$: node $(i' - 1, 2)$ cannot have an edge to node $(i', 2)$, and hence does not lie on a path from $(1, 1)$ to $(n + 1, 2)$.

Let us remark that although the main focus of real-world acronym generation systems is on the natural language parsing and interpretation of acronyms, our new algorithmic solution may inspire practical improvements in such systems or further algorithmic work.
References


A Omitted Figure

Figure 5 illustrates an example of the algorithm underlying Lemma 19.

![Diagram](image)

**Figure 5** The annotated compacted trie constructed for $T_1[i] = \{\text{abba}, \text{aaa}, \text{bb}\}$ and $T_2[j] = \{\text{ba}, \text{aaab}, \text{b}\}$ in Lemma 19. The node corresponding to $b$ has two $T_2$ labels and is an ancestor of the node corresponding to $\text{bb}$ with a $T_1$ label; hence two corresponding edges to $U_{i+1,j+1}$ are constructed. The node corresponding to $\text{aaa}$ has a $T_1$ label and is an ancestor of the node corresponding to $\text{aaab}$ with a $T_2$ label; hence a corresponding edge to $U_{i+1,j}$ is constructed. The node corresponding to $a$ has both a $T_1$ and a $T_2$ label; hence a corresponding edge to $(i+1,j+1)$ is constructed.
Compressed Indexing for Consecutive Occurrences

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Abstract

The fundamental question considered in algorithms on strings is that of indexing, that is, preprocessing a given string for specific queries. By now we have a number of efficient solutions for this problem when the queries ask for an exact occurrence of a given pattern \( P \). However, practical applications motivate the necessity of considering more complex queries, for example concerning near occurrences of two patterns. Recently, Bille et al. [CPM 2021] introduced a variant of such queries, called gapped consecutive occurrences, in which a query consists of two patterns \( P_1 \) and \( P_2 \) and a range \([a,b]\), and one must find all consecutive occurrences \((q_1, q_2)\) of \( P_1 \) and \( P_2 \) such that \( q_2 - q_1 \in [a,b] \). By their results, we cannot hope for a very efficient indexing structure for such queries, even if \( a = 0 \) is fixed (although at the same time they provided a non-trivial upper bound). Motivated by this, we focus on a text given as a straight-line program (SLP) and design an index taking space polynomial in the size of the grammar that answers such queries in time optimal up to polylog factors.

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1 Introduction

In the indexing problem, the goal is to preprocess a string for locating occurrences of a given pattern. For a string of length \( N \), structures such as the suffix tree [36] or the suffix array [31], use space linear in \( N \) and allow for answering such queries in time linear in the length of the pattern \( m \). By now, we have multiple space- and time-efficient solutions for this problem (both in theory and in practice). We refer the reader to the excellent survey by Lewenstein [29] that provides an overview of some of the approaches and some of its extensions, highlighting its connection to orthogonal range searching.
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However, from the point of view of possible applications, it is desirable to allow for more general queries than just locating an exact match of a given pattern in the preprocessed text, while keeping the time sublinear in the length of the preprocessed string. A very general query is locating a substring matching a regular expression. Very recently, Gibney and Thankachan [19] showed that if the Online Matrix-Vector multiplication conjecture holds, even with a polynomial preprocessing time we cannot answer regular expression query in sublinear time. A more reasonable and yet interesting query could concern occurrences of two given patterns that are closest to each other, or just close enough.

Preprocessing a string for queries concerning two patterns has been first studied in the context of document retrieval, where the goal is to preprocess a collection of strings. There, in the two patterns document retrieval problem the query consists of two patterns $P_1$ and $P_2$, and we must report all documents containing both of them [32]. In the forbidden pattern query problem we must report all documents containing $P_1$ but not $P_2$ [15]. For both problems, the asymptotically fastest linear-space solutions need as much as $\Omega(\sqrt{N})$ time to answer a query, where $N$ is the total length of all strings [23, 22]. That is, the complexity heavily depends on the length of the strings. Larsen et al. [28] established a connection between Boolean matrix multiplication and the two problems, thus providing a conditional explanation for the high $\Omega(\sqrt{N})$ query complexity. Later, Kopelowitz et al. [27] provided an even stronger argument using a connection to the 3SUM problem. Even more relevant to this paper is the question considered by Kopelowitz and Krauthgamer [26], who asked for preprocessing a string for computing, given two patterns $P_1$ and $P_2$, their occurrences that are closest to each other. The main result of their paper is a structure constructible in $O(N^{1.5} \log^2 N)$ time that answers such queries in $O(|P_1| + |P_2| + \sqrt{N} \log^2 N)$, for a string of length $N$, for any $\epsilon > 0$. They also established a connection between Boolean matrix multiplication and this problem, highlighting a difficulty in removing the $O(\sqrt{N})$ from both the preprocessing and query time at the same time.

The focus of this paper is the recently introduced variant of the indexing problem, called gapped indexing for consecutive occurrences, in which a query consists of two patterns $P_1$ and $P_2$ and a range $[a,b]$, and one must find the pairs of consecutive occurrences of $P_1$, $P_2$ separated by a distance in the range $[a,b]$. Navarro and Thankanchan [33] showed that for $P_1 = P_2$ there is a $O(n \log n)$-space index with optimal query time $O(m + \text{occ})$, where $m = |P_1| = |P_2|$ and occ is the number of pairs to report, but in conclusion they noticed that extending their solution to the general case of two patterns might not be possible. Bille et al. [4] provided an evidence of hardness of the general case and established a (conditional) lower bound for gapped indexing for consecutive occurrences, by connecting its complexity to that of set intersection. This lower bound suggests that, at least for indexes of size $O(N)$, achieving query time better than $\tilde{O}(|P_1| + |P_2| + \sqrt{N})$ would contradict the Set Disjointness conjecture, even if $a = 0$ is fixed. In particular, obtaining query time depending mostly on the lengths of the patterns (perhaps with some additional logarithms), arguably the whole point of string indexing, is unlikely in this case.

Motivated by the (conditional) lower bound for gapped indexing for consecutive occurrences, we consider the compressed version of this problem for query intervals $[0,b]$. For exact pattern matching, there is a long line of research devoted to designing the so-called compressed indexes, that is, indexing structures with the size being a function of the length of the compressed representation of the text, see e.g. the entry in the Encyclopedia of Algorithms [30] or the Encyclopedia of Database Systems [13]. This suggests the following research direction: can we design an efficient compressed gapped index for consecutive occurrences?
The answer of course depends on the chosen compression method. With a goal to design an index that uses very little space, we focus on the most challenging setting when the compression is capable of describing a string of exponential length (in the size of its representation). An elegant formalism for such a compression method is that of straight-line programs (SLP), which are context-free grammars describing exactly one string. SLPs are known to capture the popular Lempel-Ziv compression method up to a logarithmic factor [7, 35], and at the same time provide a more convenient interface, and in particular, allow for random access in $O(\log N)$ time [5].

By now it is known that pattern matching admits efficient indexing in SLP-compressed space. Assuming a string $S$ of length $N$ described by an SLP with $g$ productions, Claude and Navarro [9] designed an $O(g)$-space index for $S$ that allows retrieving all occurrences of a pattern of length $m$ in time $O(m^2 \log \log N + \text{occ} \log g)$. Recently, several results have improved the query time bound while still using a comparable $O(g \log N)$ amount of space: Claude, Navarro and Pacheco [10] showed an index with query time $O((m^2 + \text{occ}) \log g)$; Christiansen et al. [8] used strings attractors to further improve the time bound to $O(m + \text{occ} \log^2 N)$; and Díaz-Domínguez et al. [12] achieved $O((m \log m + \text{occ}) \log g)$ query time.

However it is not always the case that a highly compressible string is easier to preprocess. On the negative side, Abboud et al. [1] showed that, for some problems on compressed strings, such as computing the LCS, one cannot completely avoid a high dependency on the length of the uncompressed string and that for other problems on compressed strings, such as context-free grammar parsing or RNA folding, one essentially cannot hope for anything better than just decompressing the string and working with the uncompressed representation! This is also the case for some problems related to linear algebra [2]. Hence, it was not clear to us if one can avoid a high dependency on the length of the uncompressed string in the gapped indexing for consecutive occurrences problem.

In this work, we address the lower bound of Bille et al. [4] and show that, despite the negative results by Abboud et al. [1], one can circumvent it assuming that the text is very compressible:

▶ Theorem 1. For an SLP of size $g$ representing a string $S$ of length $N$, there is an $O(g^5 \log^5 N)$-space data structure that maintains the following queries: given two patterns $P_1, P_2$ both of length $O(m)$, and a range $[0, b]$, report all occ consecutive occurrences of $P_1$ and $P_2$ separated by a distance $d \in [0, b]$. The query time is $O(m \log N + (1 + \text{occ}) \cdot \log^4 N \log \log N)$.

While achieving $O(g)$ space and $O(m + \text{occ})$ query time would contradict the Set Disjointness conjecture by the reduction of Bille et al. [4], one might wonder if the space can be improved without increasing the query time and what is the true complexity of the problem when $a$ is not fixed (recall that $[a, b]$ is the range limiting the distance between co-occurrences to report). While we leave improvement on space and the general case as an interesting open question, we show that in the simpler case $a = 0, b = N$ (i.e. when there is no bound on the distance between the starting positions of $P_1$ and $P_2$), our techniques do allow for $O(g^2 \log^4 N)$ space complexity, see Corollary 11\(^1\).

Throughout the paper we assume a unit-cost RAM model of computation with word size $\Theta(\log N)$. All space complexities refer to the number of words used by a data structure.

\(^1\) Note that the conditional lower bound of Bille et al. [4] does not hold for this simpler case.
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2 Preliminaries

A string $S$ of length $|S| = N$ is a sequence $S[0]S[1] \ldots S[N-1]$ of characters from an alphabet $\Sigma$. We denote the reverse $S[N-1]S[N-2] \ldots S[0]$ of $S$ by $\text{rev}(S)$. We define $S[i \ldots j]$ to be equal to $S[i] \ldots S[j]$ which we call a substring of $S$ if $i \leq j$ and to the empty string otherwise. We also use notations $S[i \ldots j]$ and $S(i \ldots j)$ which naturally stand for $S[i] \ldots S[j-1]$ and $S[i+1] \ldots S[j]$, respectively. We call a substring $S[0 \ldots i]$ a prefix of $S$ and use a simplified notation $S[\ldots i]$, and a substring $S[i \ldots N-1]$ a suffix of $S$ denoted by $S[\ldots]$. We say that $X$ is a substring of $S$ if $X = S[i \ldots j]$ for some $0 \leq i \leq j \leq N - 1$. The index $i$ is called an occurrence of $X$ in $S$.

An occurrence $q_1$ of $P_1$ and an occurrence $q_2$ of $P_2$ form a consecutive occurrence (co-occurrence) of strings $P_1$, $P_2$ in a string $S$ if there are no occurrences of $P_1$, $P_2$ between $q_1$ and $q_2$, formally, there should be no occurrences of $P_1$ in $(q_1, q_2)$ and no occurrences of $P_2$ in $(q_1, q_2)$. For brevity, we say that a co-occurrence is $b$-close if $q_2 - q_1 \leq b$.

An integer $\pi$ is a period of a string $S$ of length $N$, if $S[i] = S[i+\pi]$ for all $i = 0, \ldots, N-1-\pi$. The smallest period of a string $S$ is called the period of $S$. We say that $S$ is periodic if the period of $S$ is at most $N/2$. We exploit the well-known corollary of the Fine and Wilf’s periodicity lemma [14]:

**Corollary 2.** If there are at least three occurrences of a string $Y$ in a string $X$, where $|X| \leq 2|Y|$, then the occurrences of $Y$ in $X$ form an arithmetic progression with a difference equal to the period of $Y$.

2.1 Grammars

**Definition 3 (Straight-line program [25]).** A straight-line program (SLP) $G$ is a context-free grammar (CFG) consisting of a set of non-terminals, a set of terminals, an initial symbol, and a set of productions, satisfying the following properties:

- A production consists of a left-hand side and a right-hand side, where the left-hand side is a non-terminal $A$ and the right-hand side is either a sequence $BC$, where $B, C$ are non-terminals, or a terminal;
- Every non-terminal is on the left-hand side of exactly one production;
- There exists a linear order $<$ on the non-terminals such that $A < B$ whenever $B$ occurs on the right-hand side of the production associated with $A$.

A run-length straight-line program (RLSLP) [34] additionally allows productions of form $A \rightarrow B^k$ for positive integers $k$, which correspond to concatenating $k$ copies of $B$. If $A$ is associated with a production $A \rightarrow a$, where $a$ is a terminal, we denote $\text{head}(A) = a$, $\text{tail}(A) = \varepsilon$ (the empty string); if $A$ is associated with a production $A \rightarrow BC$, we denote $\text{head}(A) = B$, $\text{tail}(A) = C$; and finally if $A$ is associated with a production $A \rightarrow B^k$, then $\text{head}(A) = B$, $\text{tail}(A) = B^{k-1}$.

The expansion $S$ of a sequence of terminals and non-terminals $S$ is the string that is obtained by iteratively replacing non-terminals by the right-hand sides in the respective productions, until only terminals remain. We say that $G$ represents the expansion of its initial symbol.

**Definition 4 (Parse tree).** The parse tree of a SLP (RLSLP) is a rooted tree defined as follows:
- The root is labeled by the initial symbol;
- Each internal node is labeled by a non-terminal;
- If $S$ is the expansion of the initial symbol, then the $i$th leaf of the parse tree is labeled by a terminal $S[i]$;
- A node labeled with a non-terminal $A$ that is associated with a production $A \rightarrow BC$, where $B, C$ are non-terminals, has 2 children labeled by $B$ and $C$, respectively. If $A$ is associated with a production $A \rightarrow a$, where $a$ is a terminal, then the node has one child labeled by $a$.
- (RLSLP only) A node labeled with non-terminal $A$ that is associated with a production $A \rightarrow B^k$, where $B$ is a non-terminal, has $k$ children, each labeled by $B$.

The size of a grammar is its number of productions. The height of a grammar is the

height of the parse tree. We say that a non-terminal $A$ is an ancestor of a non-terminal $B$ if there are nodes $u, v$ of the parse tree labeled with $A, B$ respectively, and $u$ is an ancestor of $v$. For a node $u$ of the parse tree, denote by $\text{off}(u)$ the number of leaves to the left of the subtree rooted at $u$.

\begin{definition}[Relevant occurrences] Let $A$ be a non-terminal associated with a production $A \rightarrow \text{head}(A)\text{tail}(A)$. We say that an occurrence $q$ of a string $P$ in $\overline{A}$ is relevant with a split $s$ if $q = |\text{head}(A)| - s \leq |\text{head}(A)| \leq q + |P| - 1$.

For example, in Fig. 1 the occurrence $q = 3$ of $P = cab$ is a relevant occurrence in $\overline{C}$ with a split $s = 1$ but $\overline{A}$ contains no relevant occurrences of $P$.
\end{definition}

\begin{claim}
Let $q$ be an occurrence of a string $P$ in a string $S$. Consider the parse tree of an RLSLP representing $S$, and let $w$ be the lowest node containing leaves $S[q], S[q+1], \ldots, S[q+|P|-1]$ in its subtree, then either
1. The label $A$ of $w$ is associated with a production $A \rightarrow BC$, and $q - \text{off}(w)$ is a relevant occurrence in $\overline{A}$; or
2. The label $A$ of $w$ is associated with a production $A \rightarrow B^r$ and $q - \text{off}(w) = q' + r' |B|$ for some $0 \leq r' \leq r$, where $q'$ is a relevant occurrence of $P$ in $\overline{A}$.

Proof. Assume first that $A$ is associated with a production $A \rightarrow BC$. We then have that the subtree rooted at the left child of $w$ (that corresponds to $B$) does not contain $S[q+|P|-1]$ and the subtree rooted at the right child of $w$ (that corresponds to $C$) does not contain $S[q]$. As a consequence, $q - \text{off}(w)$ is a relevant occurrence in $\overline{A}$.

Consider now the case where $A$ is associated with a production $A \rightarrow B^r$. The leaves labeled by $S[q]$ and $S[q+|P|-1]$ belong to the subtrees rooted at different children of $A$. If $S[q]$ belongs to the subtree rooted at the $(r'+1)$-th child of $A$, then $q' = q - \text{off}(w) - |B| \cdot r'$ is a relevant occurrence of $P$ in $\overline{A}$.
\end{claim}

\begin{definition}[Splits] Consider a non-terminal $A$ of an RLSLP $G$. If it is associated with a production $A \rightarrow BC$, define

$$\text{Splits}(A, P) = \text{Splits}_{\text{rev}}(A, P) = \{s : q \text{ is a relevant occurrence of } P \text{ in } \overline{A} \text{ with a split } s\}.$$ 

If $A$ is associated with a rule $A \rightarrow B^k$, define

$$\text{Splits}(A, P) = \{s : q \text{ is a relevant occurrence of } P \text{ in } \overline{A} \text{ with a split } s\};$$

$$\text{Splits}_{\text{rev}}(A, P) = \{|P| - s : q \text{ is a relevant occurrence of } \text{rev}(P) \text{ in } \text{rev}(\overline{A}) \text{ with split } s\}.$$ 

Define $\text{Splits}(G, P) = \text{Splits}_{\text{rev}}(G, P)$ to be the union of $\text{Splits}(A, P)$ (Splits$_{\text{rev}}(A, P)$) over all non-terminals $A$ in $G$, and $\text{Splits}'(G, P) = \text{Splits}(G, P) \cup \text{Splits}_{\text{rev}}(G, P)$.
Compressed Indexing for Consecutive Occurrences

We need the following lemma, which can be derived from Gawrychowski et al. [18]:

**Lemma 8.** Let $G$ be an SLP of size $g$ representing a string $S$ of length $N$, where $g \leq N$. There exists a Las Vegas algorithm that builds a RLSLP $G'$ of size $g' = O(g \log N)$ of height $h = O(\log N)$ representing $S$ in time $O(g \log N)$ with high probability. This RLSLP has the following additional property: For a pattern $P$ of length $m$, we can in $O(m \log N)$ time provide a certificate that $P$ does not occur in $S$, or compute the set Splits$(G', P)$. In the latter case, $|\text{Splits}'(G', P)| = O(\log N)$.

2.2 Compact Tries

We assume the reader to be familiar with the definition of a compact trie (see e.g. [21]). Informally, a trie is a tree that represents a lexicographically ordered set of strings. The edges of a trie are labeled with strings. We define the label $\lambda(u)$ of a node $u$ to be the concatenation of labels on the path from the root to $u$ and an interval $I(u)$ to be the interval of the set of strings starting with $\lambda(u)$. From the implementation point of view, we assume that a node $u$ is specified by the interval $I(u)$. The *locus* of a string $P$ is the minimum depth node $u$ such that $P$ is a prefix of $\lambda(u)$.

The standard tree-based implementation of a trie for a generic set of strings $S = \{S_1, \ldots, S_k\}$ takes $\Theta \left( \sum_{i=1}^{k} |S_i| \right)$ space. Given a pattern $P$ of length $m$ and $\tau > 0$ suffixes $Q_1, \ldots, Q_\tau$ of $P$, the trie allows retrieving the ranges of strings in (the lexicographically-sorted) $S$ prefixed by $Q_1, \ldots, Q_\tau$ in $O(m^2)$ time. However, in this work, we build the tries for very special sets of strings only, which allows for a much more efficient implementation based on the techniques of Christiansen et al. [8], the proof is given in Appendix A:

**Lemma 9.** Given an RLSLP $G$ of size $g$ and height $h$. Assume that every string in a set $S$ is either a prefix or a suffix of the expansion of a non-terminal of $G$ or its reverse. The trie for $S$ can be implemented in space $O(|S|)$ to maintain the following queries in $O(m + \tau \cdot (h + \log m))$ time: Given a pattern $P$ of length $m$ and suffixes $Q_i$ of $P$, $1 \leq i \leq \tau$, find, for each $i$, the interval of strings in (the lexicographically sorted) $S$ prefixed by $Q_i$.

3 Relevant, extremal, and predecessor occurrences in a non-terminal

In this section, we present a data structure that allows various efficient queries, which we will need to prove Theorem 1. We also show how it can be leveraged for an index in the simpler case of consecutive occurrences ($a = 0, b = N$). Recall that the text $S$ is a string of length $N$ represented by an SLP $G$ of size $g$. By applying Lemma 8, we transform $G$ into an RLSLP $G'$ of size $g' = O(g \log N)$ and depth $h = O(\log N)$ representing $S$, which we fix from now on. We start by showing that $G'$ can be processed in small space to allow multiple efficient queries:

**Theorem 10.** There is a $O(g^2 \log^4 N)$-space data structure for $G'$ that given a pattern $P$ of length $m$ can preprocess it in $O(m \log N + \log^2 N)$ time to support the following queries for a given non-terminal $A$ of $G'$:

1. Report the sorted set of relevant occurrences of $P$ in $A$ in $O(\log N)$ time;
2. Decide whether there is an occurrence of $P$ in $A$ in $O(\log N \log \log N)$ time;
3. Report the leftmost and the rightmost occurrences of $P$ in $A$, $\text{head}(A)$, and $\text{tail}(A)$ in $O(\log^2 N \log \log N)$ time;
4. Given a position $p$, find the rightmost (leftmost) occurrence $q \leq p$ ($q \geq p$) of $P$ in $A$ in $O(\log^3 N \log \log N)$ time (predecessor/successor).
Before we proceed to the proof, let us derive a data structure to report all consecutive occurrences (co-occurrences) of a given pair of patterns.

**Corollary 11.** For an SLP of size $g$ representing a string $S$ of length $N$, there is an $O(g^2 \log^4 N)$-space data structure that supports the following queries: given two patterns $P_1, P_2$ both of length $O(m)$, report all occurrences co-occurrences of $P_1$ and $P_2$ in $S$. The query time is $O(m \log N + (1 + \text{occ}) \cdot \log^3 N \log \log N)$.

**Proof.** We exploit the data structure of Theorem 10 for $G'$. To report all occurrences co-occurrences of $P_1, P_2$ in $S$, we preprocess $P_1, P_2$ in $O(m \log N + \log^2 N)$ time and then proceed as follows. Suppose that we want to find the leftmost co-occurrence of $P_1$ and $P_2$ in the string $S[i \ldots]$, where at the beginning $i = 0$. We find the leftmost occurrence $q'_1$ of $P_1$ with $q'_1 \geq t$ (if it exists) by a successor query on the initial symbol of $G'$ (the expansion of which is the entire string $S$). Then we find the leftmost occurrence $q_2$ of $P_2$ with $q_2 \geq q'_1$ (if it exists) by a successor query and the rightmost occurrence $q_1$ of $P_1$ with $q_1 \leq q_2$ by a predecessor query. If either $q'_1$ or $q_2$ do not exist, then there are no more co-occurrences in $S[i \ldots]$. Otherwise, clearly, $(q_1, q_2)$ is a co-occurrence, and there can be no other co-occurrences starting in $S[i \ldots q_2]$. In this case, we return $(q_1, q_2)$ and set $i = q_2 + 1$. The running time of the retrieval phase is $O(\log^3 N \log \log N \cdot (\text{occ} + 1))$, since we use at most three successor/predecessor queries to either output a new co-occurrence or decide that there are no more co-occurrences. ▷

### 3.1 Proof of Theorem 10

The data structure consists of two compact tries $T_{\text{pre}}$ and $T_{\text{suf}}$ defined as follows. For each non-terminal $A$, we store $\text{rev}(\text{head}(A))$ in $T_{\text{pre}}$ and $\text{tail}(A)$ in $T_{\text{suf}}$. We augment $T_{\text{pre}}$ and $T_{\text{suf}}$ by computing their heavy path decomposition:

**Definition 12.** The heavy path of a trie $T$ is the path that starts at the root of $T$ and at each node $v$ on the path branches to the child with the largest number of leaves in its subtree (heavy child), with ties broken arbitrarily. The heavy path decomposition is a set of disjoint paths defined recursively, namely it is defined to be a union of the singleton set containing the heavy path of $T$ and the heavy path decompositions of the subtrees of $T$ that hang off the heavy path.

For each non-terminal $A$ of $G'$, a heavy path $h_{\text{pre}}$ in $T_{\text{pre}}$, and a heavy path $h_{\text{suf}}$ in $T_{\text{suf}}$, we construct a multiset of points $P(A, h_{\text{pre}}, h_{\text{suf}})$. For every non-terminal $A'$ and nodes $u \in h_{\text{pre}}, v \in h_{\text{suf}}$ the multiset contains a point $(|\lambda(u)|, |\lambda(v)|)$ iff $A', u, v$ satisfy the following properties:

1. $A$ is an ancestor of $A'$;
2. $I(u)$ contains $\text{rev}(\text{head}(A'))$ and $I(v)$ contains $\text{tail}(A')$.
3. $u, v$ are the lowest nodes in $h_{\text{pre}}, h_{\text{suf}}$, respectively, satisfying Property 2.

(See Fig. 1.) The set $P(A, h_{\text{pre}}, h_{\text{suf}})$ is stored in a two-sided 2D orthogonal range emptiness data structure [29, 6] which occupies $O(|P(A, h_{\text{pre}}, h_{\text{suf}})|)$ space. Given a 2D range of the form $[\alpha, \infty] \times [\beta, \infty]$, it allows to decide whether the range contains a point in $P(A, h_{\text{pre}}, h_{\text{suf}})$ in $O(\log \log N)$ time.

**Claim 13.** The data structure occupies $O(g^2 \log^4 N)$ space.
We then sort the Splits of \( u \) for \( T_{\text{pre}} \) and \( T_{\text{suf}} \) respectively. The heavy paths \( h_{\text{pre}} \) in \( T_{\text{pre}} \) and \( h_{\text{suf}} \) in \( T_{\text{suf}} \) are shown in blue. We have \((2,2) \in \mathcal{P}(A, h_{\text{pre}}, h_{\text{suf}})\) corresponding to \( C, u, v \).

**Proof.** Each non-terminal \( A' \) has at most \( g' \) distinct ancestors and each root-to-leaf path in \( T_{\text{pre}} \) or \( T_{\text{suf}} \) crosses \( O(\log g') \) heavy paths (as each time we switch heavy paths, the number of leaves in the subtree of the current node decreases by at least a factor of two). As a corollary, each non-terminal creates \( O(g \log^2 g') = O(g \log^3 N) \) points across all orthogonal range emptiness data structures.

When we receive a pattern \( P \), we compute \( \text{Splits}'(G', P) \) via Lemma 8 in \( O(m \log N) \) time or provide a certificate that \( P \) does not occur in \( S \), in which case there are no occurrences of \( P \) in the expansions of the non-terminals of \( G' \). Recall that \( |\text{Splits}'(G', P)| \in O(\log N) \). We then sort \( \text{Splits}'(G', P) \) in \( O(\log^2 N) \) time (a technicality which will allow us reporting relevant occurrences sorted without time overhead). Finally, we compute, for each \( s \in \text{Splits}'(G', P) \), the interval of strings in \( T_{\text{pre}} \) prefixed by \( \text{rev}(P[\ldots s]) \) (which is the interval \( I(u) \) for the locus \( u \) of \( \text{rev}(P[\ldots s]) \) in \( T_{\text{pre}} \)) and the interval of strings in \( T_{\text{suf}} \) prefixed by \( P(s \ldots) \) (which is the interval \( I(u) \) for the locus \( u \) of \( P(s \ldots) \) in \( T_{\text{suf}} \)). By Lemma 9, with \( \tau = |\text{Splits}'(G', P)| = O(\log N) \) and \( h = O(\log N) \), this step takes \( O(m + \log^2 N) \) time.

Reporting relevant occurrences is easy: by definition, each relevant occurrence \( q \) of \( P \) in \( \mathcal{A} \) is equal to \( [\text{head}(A)] - s \) for some \( s \in \text{Splits}'(G', P) \) such that \( \text{rev}(P[\ldots s]) \) is a prefix of \( \text{rev}(\text{head}(A)) \) and \( P(s \ldots) \) is a prefix of \( \text{tail}(A) \). As we already know the intervals of the strings in \( T_{\text{suf}} \) and \( T_{\text{pre}} \), starting with \( \text{rev}(P[\ldots s]) \) and \( P(s \ldots) \), respectively, both conditions can be checked in constant time per split, or in \( O(|\text{Splits}'(G', P)|) = O(\log N) \) time overall. Note that since \( \text{Splits}'(G', P) \) are sorted, the relevant occurrences are reported sorted as well.

We now explain how to answer emptiness queries on a non-terminal:

\begin{itemize}
  \item \textbf{Claim 14.} Let \( A \) be a non-terminal labeling a node in the parse tree of \( G' \). We can decide whether \( \mathcal{A} \) contains an occurrence of \( P \) in \( O(\log N \log \log N) \) time.
\end{itemize}

**Proof.** Below we show that \( P \) occurs in \( \mathcal{A} \) if and only if there exists a split \( s \in \text{Splits}'(G', P) \) such that for \( u \) being the locus of \( \text{rev}(P[\ldots s]) \) in \( T_{\text{pre}} \) and \( v \) the locus of \( P(s \ldots) \) in \( T_{\text{suf}} \), for \( h_{\text{pre}} \) the heavy path containing \( u \) in \( T_{\text{pre}} \) and \( h_{\text{suf}} \) the heavy path containing \( v \) in \( T_{\text{suf}} \), the rectangle \([|\lambda(u)|, +\infty] \times [|\lambda(v)|, +\infty] \) contains a point from \( \mathcal{P}(A, h_{\text{pre}}, h_{\text{suf}}) \). Before we proceed to the proof, observe that by the bound on \( |\text{Splits}'(G', P)| \) this allows us to decide whether \( P \) occurs in \( \mathcal{A} \) in \( O(\log N) \) range emptiness queries, which results in \( O(\log N \log \log N) \) query time.

Assume that \([|\lambda(u)|, +\infty] \times [|\lambda(v)|, +\infty] \) contains a point \((x, y) \in \mathcal{P}(A, h_{\text{pre}}, h_{\text{suf}})\) corresponding to a non-terminal \( A' \). By construction, \( A \) is an ancestor of \( A' \), the subtree of \( u \) contains a leaf corresponding to \( \text{rev}(\text{head}(A')) \) and the subtree of \( v \) contains a leaf corresponding to \( \text{tail}(A') \). Consequently, \( \mathcal{A} \) contains an occurrence of \( P \), which implies...
that $\overline{A}$ contains an occurrence of $P$. To show the reverse direction, let $\ell = \text{off}(u) + 1$ and $r = \text{off}(u) + |\overline{A}|$, i.e. $S[\ell \ldots r] = \overline{A}$. The string $\overline{A}$ contains an occurrence $\overline{A}[q \ldots q + |P|]$ of $P$ iff $S[\ell + q \ldots \ell + q + |P|]$ is an occurrence of $P$ in $S$. From Claim 6 it follows that if $w$ is the lowest node in the parse tree of $G'$ that contains leaves $S[\ell + q], \ldots, S[\ell + q + |P| - 1]$ in its subtree and $A'$ is its label, then there exists a split $s \in \text{Splits}'(G', P)$ such that $\text{rev}(P[\ldots s])$ is a prefix of $\text{rev}(\text{head}(A'))$ and $P(s \ldots]$ of $\text{tail}(A')$. By definition of $u$ and $v$, the leaf of $T_{pre}$ labeled with $\text{rev}(\text{head}(A'))$ belongs to $I(u)$ and the leaf of $T_{suf}$ labeled with $\text{tail}(A')$ belongs to $I(v)$. Let $h_{pre}$ $(h_{suf})$ be the heavy path in $T_{pre}(T_{suf})$ containing $u$ $(v)$ and $(x, y)$ be the point in $\mathcal{P}(A, h_{pre}, h_{suf})$ created for $A'$. As $|\lambda(u)| \leq x$ and $|\lambda(v)| \leq y$, the rectangle $[[\lambda(u)], +\infty] \times [[\lambda(v)], +\infty]$ is not empty. \hfill $<$

It remains to explain how to retrieve the leftmost/rightmost occurrences in a non-terminal, as well as to answer predecessor/successor queries. The main idea for all four types of queries is to start at any node of the parse tree of $G'$ labeled by $A$ and recurse down via emptiness queries and case inspection. Since the length of the expansion decreases each time we recurse from a non-terminal to its child and the height of $G'$ is $h = O(\log N)$, this allows to achieve the desired query time. We provide full details in Appendix B.

4 Compressed Indexing for Close Co-occurrences

In this section, we show our main result, Theorem 1. Recall that $S$ is a string of length $N$ represented by an SLR $G$ of size $g$. We start by applying Lemma 8 to transform $G$ into an RLSLP $G'$ of size $g' = O(g \log N)$ and height $h = O(\log N)$ representing $S$.

The query algorithm uses the following strategy: first, it identifies all non-terminals of $G'$ such that their expansion contains a $b$-close relevant co-occurrence, where a relevant co-occurrence is defined similarly to a relevant occurrence:

\begin{definition} [Relevant co-occurrence] Let $A$ be a non-terminal of $G'$. We say that a co-occurrence $(q_1, q_2)$ of $P_1, P_2$ in $\overline{A}$ is relevant if $q_1 \leq |\text{head}(A)| \leq q_2 + |P_2| - 1$.
\end{definition}

Second, it retrieves all $b$-close relevant co-occurrences in each of those non-terminals, and finally, reports all $b$-close co-occurrences by traversing the (pruned) parse tree of $G'$, which is possible due to the following claim:

\begin{claim} Assume that $P_2$ is not a substring of $P_1$, and let $(q_1, q_2)$ be a co-occurrence of $P_1, P_2$ in a string $S$. In the parse tree of $G'$, there exists a unique node $u$ such that either

1. Its label $A$ is associated with a production $A \rightarrow BC$, and $(q_1 - \text{off}(u), q_2 - \text{off}(u))$ is a relevant co-occurrence of $P_1, P_2$ in $\overline{A}$;

2. Its label $A$ is associated with a production $A \rightarrow B^k$, $q_1 - \text{off}(u) = q'_1 + k'|\overline{B}|$, $q_2 - \text{off}(u) = q'_2 + k'|\overline{B}|$ for some $0 \leq k' \leq k$, where $(q'_1, q'_2)$ is a relevant co-occurrence of $P_1, P_2$ in $\overline{A}$.
\end{claim}

Proof. Let $A$ be the label of the lowest node $u$ in the parse tree that contains leaves $S[q_1], S[q_1 + 1], \ldots, S[q_2 + |P_2| - 1]$ in its subtree. Because $P_2$ is not a substring of $P_1$, $A$ cannot be associated with a production $A \rightarrow a$. By definition, $S[\text{off}(u) + 1]$ is the leftmost leaf in the subtree of this node.

Assume first that $A$ is associated with a production $A \rightarrow BC$. We then have that the subtree rooted at the left child of $u$ (labeled by $B$) does not contain $S[q_2 + |P_2| - 1]$ and the subtree rooted at the right child of $u$ (labeled by $C$) does not contain $S[q_1]$. As a consequence, $(q_1 - \text{off}(u), q_2 - \text{off}(u))$ is a relevant co-occurrence of $P_1, P_2$ in $\overline{A}$.\hfill $<$

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Consider now the case where $A$ is associated with a production $A \rightarrow B^k$. The leaves labeled by $S[q_1]$ and $S[q_2 + |P_2| - 1]$ belong to the subtrees rooted at different children of $A$. If $S[q_1]$ belongs to the subtree rooted at the $k$-th child of $A$, then $(q_1 - \text{off}(u) - |\overline{B}| \cdot (k' - 1), q_2 - \text{off}(u) - |\overline{B}| \cdot (k' - 1))$ is a relevant co-occurrence of $P_1, P_2$ in $\overline{A}$.

### 4.1 Combinatorial observations

Informally, we define a set of $O(g^2)$ strings and show that for any patterns $P_1, P_2$ there are two strings $S_1, S_2$ in the set with the following property: whenever the expansion of a non-terminal $A$ in $G'$ contains a pair of occurrences $P_1, P_2$ forming a relevant co-occurrence, there are occurrences of $S_1, S_2$ in the proximity. This will allow us to preprocess the non-terminals of $G'$ for occurrences of the strings in the set and use them to detect $b$-close relevant co-occurrences of $P_1, P_2$.

Consider two tries, $T_{\text{pre}}$ and $T_{\text{suf}}$: For each production of $G'$ of the form $A \rightarrow BC$, we store $\overline{C}$ in $T_{\text{suf}}$ and $\text{rev}(\overline{B})$ in $T_{\text{pre}}$. For each production of the form $A \rightarrow B^k$, we store $\overline{B}$, $\overline{B^2}$, $\overline{B^{k-2}}$, and $\overline{B^{k-1}}$ in $T_{\text{suf}}$ and the reverses of those strings in $T_{\text{pre}}$. For $j \in \{1, 2\}$ and $s \in \text{Splits}'(G', P_j)$ define $S_j(s) = \text{rev}(U)V$, where $U$ is the label of the locus of $\text{rev}(P_j[s\ldots])$ in $T_{\text{pre}}$ and $V$ is the label of the locus of $P_j(s\ldots)$ in $T_{\text{suf}}$. Let $l_j(s) = |\text{rev}(U)|$ and $\Delta_j(s) = l_j(s) - s$.

Consider a non-terminal $A$ such that its expansion $A$ contains a relevant co-occurrence $(q_1, q_2)$ of $P_1, P_2$.

\textbf{Claim 17.} There exists $s \in \text{Splits}'(G', P_2)$ such that $p_2 = q_2 - \Delta_2(s)$ is an occurrence of $S_2(s)$ in $\overline{A}$ and $|p_2, p_2 + |S_2(s)|| \geq |q_2, q_2 + |P_2||$.

\textbf{Proof.} Below we show that there exists a descendant $A'$ of $A$ and a split $s \in \text{Splits}'(G', P_2)$ such that either $\text{rev}(P_2[s\ldots])$ is a prefix of $\text{rev}(\text{head}(A'))$ and $P_2(s\ldots)$ is a prefix of $\text{tail}(A')$, or $A'$ is associated with a rule $A' \rightarrow (B')^k$, $\text{rev}(P_2[s\ldots])$ is a prefix of $\text{rev}((B')^2)$ and $P_2(s\ldots)$ is a prefix of $(B')^{k-2}$. The claim follows by the definition of $T_{\text{pre}}, T_{\text{suf}}$, and $S_2(s)$.

If $q_2$ is relevant in $\overline{A}$, there exists a split $s \in \text{Splits}'(G', P_2)$ such that $\text{rev}(P_2[s\ldots])$ is a prefix of $\text{rev}(\text{head}(A))$ and $P_2(s\ldots)$ is a prefix of $\text{tail}(A)$ by definition. If $q_2$ is not relevant, then $q_2 \geq |\text{head}(A)|$ by the definition of a co-occurrence. By Claim 6, there is a descendant $A'$ of $A$ corresponding to a substring $A[\ell\ldots r]$ for which either $(q_2 - \ell)$ is relevant (and then we can repeat the argument above), or $A'$ is associated with a rule $A' \rightarrow (B')^k$ and $(q_2 - \ell) - k' \cdot |\overline{B^2}|$ is relevant, for some $0 \leq k' \leq k$. Consider the latter case. If $A' = A$, then $k' = 1$, as otherwise $q_1 < q_2' = q_2 - |\overline{B^2}| < q_3$ is an occurrence of $P_2$ in $\overline{A}$ contradicting the definition of a co-occurrence (recall that $(q_1, q_2)$ is a relevant co-occurrence and hence by definition $q_1 < |\text{head}(A)|$), and therefore $s = |B'|^2 - q_2 + \ell \in \text{Splits}'(G', P_2)$, $\text{rev}(P_2[s\ldots])$ is a prefix of $\text{rev}(B'^2)$ and $P_2(s\ldots)$ is a prefix of $(B')^{k-2}$. If $A' \neq A$, then we can analogously conclude that $k' = 0$, which implies $s = |\overline{B^2}| - q_2 + \ell \in \text{Splits}'(G', P_2)$, $\text{rev}(P_2[s\ldots])$ is a prefix of $\text{rev}(B')$ and $P_2(s\ldots)$ is a prefix of $(B')^{k-1}$.

As the definition of a co-occurrence is not symmetric, $q_1$ does not enjoy the same property. However, a similar claim can be shown:

\textbf{Lemma 18.} There exists $s \in \text{Splits}'(G', P_1)$ and an occurrence $p_1$ of $S_1(s)$ in $\overline{A}$ such that $|p_1, p_1 + |S_1(s)|| \geq |q_1, q_1 + |P_1||$ and at least one of the following holds:

1. $q_1 - \Delta_1(s)$ is an occurrence of $S_1(s)$;
2. $q_2$ is a relevant occurrence of $P_2$ in $\overline{A}$, the period of $S_1(s)$ equals the period $\pi_1$ of $P_1$, and there exists an integer $k$ such that $p_1 = q_1 - \Delta_1(s) - \pi_1 \cdot k$ and $q_2 + \pi_1 - 1 \leq p_1 + |S_1(s)| - 1 \leq q_2 + |P_2| - 1$. 


Proof. If \( q_1 \) is a relevant occurrence of \( P_1 \) in \( A \) with a split \( s \in \text{Splits}'(G', P_1) \), then \( \text{rev}(P_1[\ldots s]) \) is a prefix of \( \text{rev}(\text{head}(A)) \) and \( P_1(s \ldots] \) is a prefix of \( \text{tail}(A) \) and therefore the first case holds by the definition of \( T_{\text{pre}} \) and \( T_{\text{surf}} \).

Otherwise, by Claim 6, there is a descendant \( A' \) of \( \text{head}(A) \) corresponding to a substring \( \overline{A}[\ell \ldots r] \) for which either \( (q_1 - \ell) \) is relevant (and then we can repeat the argument above), or \( A' \) is associated with a rule \( A' \rightarrow (B')^k \) and \( (q_1 - \ell) - k' \cdot \lvert B' \rvert \), for some \( 0 \leq k' \leq k \), is a relevant occurrence of \( P_1 \) in \( A' \) with a split \( s \in \text{Splits}'(G', P_1) \). Consider the latter case. We must have (1) \( q_1 + |P_1| - 1 + |B'| \geq r \) or (2) \( q_1 + |B'| - 1 \geq q_2 \), because if both inequalities do not hold, then \( q_1 < q_1 + |B'| \leq q_2 \) is an occurrence of \( P_1 \) in \( A \), which contradicts the definition of a co-occurrence. Additionally, if (1) holds, then by definition there exists a split \( s' \in \text{Splits}'(G', P_1) \) (which might be different from the split \( s \) above) such that \( \text{rev}(P_1[\ldots s']) \) is a prefix of \( \text{rev}(B')^{r-1} \) and \( P_1(s' \ldots] \) is a prefix of \( B' \) and we fall into the first case of the lemma.

From now on, assume that (2) holds and (1) does not. Since \( q_1 + |B'| \leq r \leq \lvert \text{head}(A) \rvert \) and \( (q_1, q_2) \) is a relevant co-occurrence, \( q_2 \) must be a relevant occurrence of \( P_2 \) in \( A' \). If \( |P_1| - s \leq \lvert (B')^2 \rvert \), then \( \text{rev}(P_1[\ldots s]) \) is a prefix of \( \text{rev}(B')^2 \) and \( P_1(s \ldots] \) is a prefix of \( (B')^2 \). Furthermore, \( q_1 = q_1 - \Delta_1(s) \) is an occurrence of \( S_1(s) \). Otherwise, by Fine and Wilf's periodicity lemma [14], the periods of \( A', P_1, \) and \( S_1(s) \) are equal, since \( P_1 \) and hence \( S_1(s) \) span at least two periods of \( A' \). By periodicity, \( S_1(s) \) occurs at positions \( q_1 - \Delta_1(s) - |B'| \cdot k \) of \( A' \). Let \( p_1 \) be the leftmost of these positions which satisfies \( p_1 + |S_1(s)| - 1 \geq q_1 + |P_1| - 1 \). This position is well-defined as (1) does not hold, and furthermore \( [q_1, q_1 + |P_1|] \subseteq [p_1, p_1 + |S_1(s)|] \) as \( s \leq t_1(s) \) and \( |S_1(s)| - t_1(s) \geq |P_1| - s \). We have \( p_1 = q_1 - \Delta_1(s) - \pi, k \) for some integer \( k \) (as \( |B'| \) is a multiple of \( \pi_1 \)), and \( q_2 + \pi_1 - 1 \leq q_2 + |B'| - 1 \leq q_2 + |P_1| - 1 \leq p_1 + |S_1(s)| - 1 \leq r < q_2 + |P_2| - 1 \), where the last inequality holds as \( q_2 \) is a relevant occurrence in \( A' \). The claim of the lemma follows. 

We summarize Claim 17 and Lemma 18:

**Corollary 19.** Let \((q_1, q_2)\) be a co-occurrence of \( P_1, P_2 \) in the expansion of a non-terminal \( A \). There exist splits \( s_1 \in \text{Splits}'(G', P_1), s_2 \in \text{Splits}'(G', P_2) \) and occurrences \( p_1 \) of \( S_1(s_1) \) and \( p_2 \) of \( S_2(s_2) \), where \([p_1, p_1 + |S_1(s_1)|] \supseteq [q_1, q_1 + |P_1|] \) and \([p_2, p_2 + |S_2(s_2)|] \supseteq [q_2, q_2 + |P_2|], \) such that at least one of the following holds:

1. The occurrence \( p_1 \) is either relevant or \( p_1 + |S_1(s_1)| - 1 \leq \lvert \text{head}(A) \rvert \). The occurrence \( p_2 \) is either relevant or \( p_2 > \lvert \text{head}(A) \rvert \). Additionally, \( p_1 = q_1 - \Delta_1(s_1) \) and \( p_2 = q_2 - \Delta_2(s_2) \).

2. The occurrence \( p_2 \) is relevant and \( p_1 \leq \lvert \text{head}(A) \rvert \). Additionally, \( p_2 = q_2 - \Delta_2(s_2) \), the period of \( S_1(s) \) equals the period \( \pi_1 \) of \( P_1 \), and there exists an integer \( k \) such that \( p_1 = q_1 - \Delta_1(s_1) - \pi_1 \cdot k \) and \( p_2 + \pi_1 - 1 \leq p_1 + |S_1(s_1)| - 1 \leq p_2 + |S_2(s_2)| - 1 \).

The reverse observation holds as well:

**Observation 20.** If \( p_j \) is an occurrence of \( S_j(s_j) \) in \( A \), \( j = 1, 2 \), then \( q_j = p_j + \Delta_j(s_j) \) is an occurrence of \( P_j \). Furthermore, if \( S_1(s) \) is periodic with period \( \pi_1 \), then \( q_1 + \pi_1 \cdot k, 0 \leq k \leq [\lvert S_1(s) \rvert - q_1 - |P_1|] / \pi_1 \), are occurrences of \( P_1 \) in \( A \).

Finally, the following trivial observation will be important for upper bounding the time complexity of our query algorithm:

**Observation 21.** If a string contains a pair of occurrences \((q_1, q_2)\) of \( P_1 \) and \( P_2 \) such that \( 0 \leq q_2 - q_1 \leq b \), then it contains a \( b \)-close co-occurrence of \( P_1 \) and \( P_2 \).
Compressed Indexing for Consecutive Occurrences

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(a) If neither (1) nor (2) then \((q_1, q_2)\) is not consecutive.

(b) (1) holds and (2) does not.

(c) (2) holds, (1) does not, and \(|P_1| - s \geq (B')^2\).

\[\text{Figure 2 Subcases of Lemma 18.}\]

4.2 Index

The first part of the index is the data structure of Theorem 10 and the index of Christiansen et al. [8]:

\[\text{Fact 22 ([8, Introduction and Theorem 6.12]).} \text{ There is a } O(q \log^2 N)\text{-space data structure that can find the occ occurrences of any pattern } P[1 \ldots m] \text{ in } S \text{ in time } O(m + occ).}\]

The second part of the index are the tries \(T_{pre}\) and \(T_{suf}\), augmented as explained below. Consider a quadruple \((u_1, u_2, v_1, v_2)\), where \(u_1\) and \(u_2\) are nodes of \(T_{pre}\) and \(v_1\) and \(v_2\) are nodes of \(T_{suf}\). Let \(U_1, U_2, V_1, V_2\) be the labels of \(u_1, u_2, v_1, v_2\), respectively. Define \(S_1 = \text{rev}(U_1)V_1\) and \(S_2 = \text{rev}(U_2)V_2\), and let \(l_1 = |\text{rev}(U_1)|\) and \(l_2 = |\text{rev}(U_2)|\).

First, we store a binary search tree \(T_l(u_1, u_2, v_1, v_2)\) that for each non-terminal \(A\) contains at most six integers \(d = p_2 - p_1\), where \(p_1, p_2\) are occurrences of \(S_1, S_2\) in \(A\), satisfying at least one of the below:

1. \(p_1\) is the rightmost occurrence of \(S_1\) such that \(p_1 + |S_1| - 1 < |\text{head}(A)|\) and \(p_2\) is the leftmost occurrence of \(S_2\) such that \(p_2 \geq |\text{head}(A)|\);
2. \(p_1\) is a relevant occurrence of \(S_1\) with a split \(l_1\) and \(p_2\) is the leftmost occurrence of \(S_2\) such that \(p_2 \geq |\text{head}(A)|\);
3. \(p_1\) is a relevant occurrence of \(S_1\) with a split \(l_1\), \(p_2\) is a relevant occurrence of \(S_2\) with a split \(l_2\);
4. \(p_2\) is a relevant occurrence of \(S_2\) with a split \(l_2\) and \(p_1\) is the rightmost occurrence of \(S_1\) such that \(p_1 + |S_1| - 1 < p_2\);
5. \(p_2\) is a relevant occurrence of \(S_2\) with a split \(l_2\) and \(p_1\) is the leftmost or second leftmost occurrence of \(S_1\) in \(\text{head}(A)\) such that \(p_1 < p_2 \leq p_1 + |S_1| - 1 < p_2 + |S_2| - 1\).
Second, we store a list of non-terminals $L(u_2, v_2)$ such that their expansion contains a relevant occurrence of $S_2$ with a split $l_2$. Additionally, for every $k \in [0, \log N]$, we store, if defined:
1. The rightmost occurrence $p_1$ of $S_1$ in $S_2$ such that $p_1 + (|S_1| - 1) \leq l_2 - 2^k$;
2. The leftmost occurrence $p'_1$ of $S_1$ in $S_2$ such that $p'_1 \leq l_2 - 2^k \leq p_1 + |S_1| - 1$;
3. The rightmost occurrence $p''_1$ of $S_1$ in $S_2$ such that $p''_1 \leq l_2 - 2^k \leq p'_1 + |S_1| - 1$.

Finally, we compute and memorize the period $\pi_1$ of $S_1$. If the period is well-defined (i.e., $S_1$ is periodic), we build a binary search tree $T_2(u_1, u_2, v_1, v_2)$. Consider a non-terminal $A$ containing a relevant occurrence $p_2$ of $S_2$ with a split $l_2$. Let $p_1$ be the leftmost occurrence of $S_1$ such that $p_1 \leq p_2 \leq p_1 + |S_1| - 1 \leq p_2 + |S_2| - 1$ and $p'_1$ the rightmost. If $p_1$ and $p'_1$ exist ($p_1$ might be equal to $p'_1$) and $p'_1 + |S_1| - 1 \geq p_2 + \pi_1 - 1$, we add an integer $(p'_1 - p_1)/\pi_1$ to the tree and associate it with $A$. We also memorize a number $ov(S_1, S_2) = p_2 - p'_1$, which does not depend on $A$ by Corollary 2 and therefore is well-defined (it corresponds to the longest prefix of $S_2$ periodic with period $\pi_1$).

> Claim 23. The data structure occupies $O(g^5 \log^5 N)$ space.

Proof. The data structure of Theorem 10 occupies $O(g^2 \log^4 N)$ space. The index of Christiansen et al. occupies $O(g \log^2 N)$ space. There are $O((g')^4)$ quadruples $(u_1, u_2, v_1, v_2)$ and for each of them the trees take $O(g')$ space. The arrays of occurrences of $S_1$ in $S_2$ use $O(\log N)$ space. Therefore, overall the data structure uses $O(g^5 \log^5 N)$ space. \hfill ◼

4.3 Query

Recall that a query consists of two strings $P_1, P_2$ of length at most $m$ each and an integer $b$, and we must find all $b$-close co-occurrences of $P_1, P_2$ in $S$, let occ be their number.

We start by checking whether $P_2$ occurs in $P_1$ using a linear-time and constant-space pattern matching algorithm such as [11]. If it is, let $q_2$ be the position of the first occurrence. If $q_2 > b$, then there are no $b$-close co-occurrences of $P_1, P_2$ in $S$. Otherwise, to find all $b$-close co-occurrences of $P_1, P_2$ in $S$ (that always consist of an occurrence of $P_1$ in $S$ and the first occurrence of $P_2$ in $P_1$), it suffices to find all occurrences of $P_1$ in $S$, which we do using the index of Christiansen et al. [8] in time $O(|P_1| + \text{occ}) = O(m + \text{occ})$.

From now on, assume that $P_2$ is not a substring of $P_1$. Let $\mathcal{N}'$ be the set of all non-terminals in $G'$ such that their expansion contains a relevant $b$-close co-occurrence of $P_1, P_2$. By Claim 16, $|\mathcal{N}'| \leq \text{occ}$.

> Lemma 24. Assume that $P_2$ is not a substring of $P_1$. One can retrieve in $O(m + (1 + \text{occ}) \log^3 N)$ time a set $\mathcal{N}' \supset \mathcal{N}'$, $|\mathcal{N}'| = O(\text{occ} \log N)$.

Proof. We start by computing $\text{Splits}'(G', P_1)$ and $\text{Splits}'(G', P_2)$ via Lemma 8 in $O(|P_1| + |P_2| \log N) = O(m \log N)$ time (or providing a certificate that either $P_1$ or $P_2$ does not occur in $S$, in which case there are no co-occurrences of $P_1, P_2$ in $S$ and we are done). Recall that $|\text{Splits}'(G', P_1)|, |\text{Splits}'(G', P_2)| \in O(\log N)$. For each fixed pair of splits $s_1 \in \text{Splits}'(G', P_1)$, $s_2 \in \text{Splits}'(G', P_2)$ and $j \in [1, 2]$, we compute the interval of strings in $T_{\text{pre}}$ prefixed by rev$(P_j[s_1, \ldots, s_j])$, which corresponds to the locus $u_j = \text{rev}(P_j[s_1, \ldots, s_j])$ in $T_{\text{pre}}$ and the interval of strings in $T_{\text{post}}$ prefixed by $P_j(s_j \ldots)$, which corresponds to the locus $v_j = P_j(s_j \ldots)$ in $T_{\text{post}}$. Computing the intervals takes $O(m + \log^2 N)$ time for all the splits by Lemma 9.
Consider the strings $S_1 = \text{rev}(U_1)V_1$ and $S_2 = \text{rev}(U_2)V_2$, where $U_1, U_2, V_1, V_2$ are the labels of $u_1, v_1, u_2, v_2$, respectively. Let $l_1 = |\text{rev}(U_1)|$, $\Delta_1 = l_1 - s_1$, $l_2 = |\text{rev}(U_2)|$, $\Delta_2 = l_2 - s_2$, and $\Delta = \Delta_1 - \Delta_2$.

Consider a relevant co-occurrence $(q_1, q_2)$ of $P_1, P_2$ in the expansion of a non-terminal $A$. By Corollary 19, $q_1, q_2$ imply existence of occurrences $p_1, p_2$ of $S_1, S_2$ such that $|p_1| + |S_1| \geq |q_1| + |P_1|$ and $|p_2| + |S_2| \geq |q_2| + |P_2|$. Our index must treat both cases of Corollary 19. We consider eight subcases defined in Fig. 3, which describe all possible locations of $p_1$ and $p_2$.

**Figure 3** Assume that $S_1$ does not contain $S_2$. The figure shows all possible locations of occurrences $p_1, p_2$ of $S_1, S_2$ in $A$. In Case 1 of Corollary 19, there are six subcases: (1.1) $p_1 + |S_1| - 1 \leq |\text{head}(A)|$, $p_2 > |\text{head}(A)|$; (1.2) $p_1$ is a relevant occurrence of $S_1$, $p_2 > |\text{head}(A)|$; (1.3) $p_1, p_2$ are relevant; (1.4) $p_2$ is relevant, $p_1 + |S_1| - 1 \leq p_2$; (1.5) $p_2$ is relevant, $p_2 < p_1 \leq p_1 + |S_1| - 1 \leq p_2 + |S_2| - 1$; (1.6) $p_2$ is relevant, $p_1 < p_2 < p_1 + |S_1| - 1 \leq p_2 + |S_2| - 1$. By the definition of a co-occurrence and by Observation 20, in Subcases (1.1) and (1.4) $p_1$ must be as far to the right as possible, and in Subcases (1.1) and (1.2) $p_2$ must be as far to the left as possible. In Case 2, there are two subcases: (2.1) $p_2$ is relevant and $p_2 \leq p_1 \leq p_1 + |S_1| - 1 \leq p_2 + |S_2| - 1$; (2.2) $p_2$ is relevant and $p_1 < p_2 < p_1 + \pi_1 - 1 \leq p_1 + |S_1| - 1$, where $\pi_1$ is the period of $S_1$. In all subcases, $q_2 = p_2 + \Delta_2$. In Subcases (1.1)-(1.6) $q_1 = p_1 + \Delta_1$ and in Subcases (2.1) and (2.2) $q_1 = p_1 + \Delta_1 + k \cdot \pi_1$ for some integer $k$.

**Subcases (1.1)-(1.4).** To retrieve the non-terminals, we query $T_1(u_1, v_1, u_1, v_1)$ to find all integers that belong to the range $[\Delta, \Delta + b]$ (and the corresponding non-terminals). Recall that, for each non-terminal $A$, the tree stores an integer $d = p_2 - p_1$, where $p_1$ is the starting position of an occurrence of $S_1$ in $A$ and $p_2$ of $S_2$. By Observation 20, $p_1 + \Delta_1$ is an occurrence of $P_1$ and $p_2 + \Delta_2$ is an occurrence of $P_2$. The distance between them is in $[0, b]$ if $d \in [\Delta, \Delta + b]$. By Observation 21, each retrieved non-terminal contains a close co-occurrence of $(q_1, q_2)$. On other hand, if $A$ contains a co-occurrence $(q_1, q_2)$ corresponding to one Subcases (1.1)-(1.4), then by Corollary 19, $p_1 = q_1 - \Delta_1$ is an occurrence of $S_1$ and $p_2 = q_2 - \Delta_2$ is an occurrence of $S_2$ and by construction $T_1(u_1, v_2, v_1, v_2)$ stores an integer $d = p_2 - p_1$. Therefore, the query retrieves all non-terminals corresponding to Subcases (1.1)-(1.4).
Subcases (1.5) and (2.1). We must decide whether an occurrence of $P_1$ in $S_2$ forms a $b$-close co-occurrence with the occurrence $\Delta_2$ of $P_2$ in $S_2$, and if so, report all non-terminals such that their expansion contains a relevant co-occurrence of $S_2$ with a split $l_2$, which are exactly the non-terminals stored in the list $L(u_2,v_2)$. Let $k = \lfloor \log(s_2) \rfloor$. Recall that the index stores the following information for $k$:

1. $p_1$, the rightmost occurrence of $S_1$ in $S_2$ such that $p_1 + (|S_1| - 1) \leq l_2 - 2^k$;
2. $p'_1$, the leftmost occurrence of $S_1$ in $S_2$ such that $p'_1 \leq l_2 - 2^k \leq p_1 + (|S_1| - 1)$;
3. $p''_1$, the rightmost occurrence of $S_1$ in $S_2$ such that $p''_1 \leq l_2 - 2^k \leq p_1 + (|S_1| - 1)$.

(See Fig. 4). By Observation 20, the occurrence $p_1$ of $S_1$ induces an occurrence $q_1 = p_1 + \Delta_1$ of $P_1$. Furthermore, if $S_1$ is periodic with period $\pi_1$, then $q_1 + \pi_1 \leq 0 \leq [([|S_1| - q_1 - |P_1|]/\pi_1)$, are also occurrences of $P_1$. One can decide whether the distance from any of these occurrences to $q_2$ is in $[0,b]$ in constant time, and if yes, then there $S_2$ contains a $b$-close co-occurrence of $P_1,P_2$ by Observation 21. Second, by Corollary 2, if $S_1$ is not periodic, then there are no occurrences of $S_1$ between $p'_1$ and $p''_1$ such that $p''_1$ by Observation 20 induce occurrences $p'_1 + \Delta_1, p''_1 + \Delta_1$ of $P_1$. Otherwise, there are occurrences of $P_1$ in every position $p'_1 + \Delta_1 + k \cdot \pi_1$, $0 \leq k \leq ([([|S_1| + p''_1 - |P_1| - p'_1]/\pi_1$). Similarly, we can decide whether the distance from any of them to the occurrence $\Delta_2$ of $P_2$ in $S_2$ is in $[0,b]$ in constant time. Finally, let $q_1$ be the rightmost occurrence of $P_1$ in $S_2$ in the interval $[l_2 - 2^k + 1, \Delta_2]$. We extract $S_2(l_2 - 2^k, \Delta_2 + |P_2|)$ via Fact 28 and search for $q_1$ using a linear-time pattern matching algorithm for $P_1$, which takes $O(|P_1| + |P_2|) = O(m)$ time. If $0 \leq \Delta_2 - q_1 \leq b$, then there is a $b$-close co-occurrence of $P_1,P_2$ in $S_2$. Correctness follows from Corollary 19, Observation 20 and Observation 21.

**Figure 4** Query algorithm for Subcases (1.5) and (2.1).

Subcase (2.2). Let $\pi_1$ be the period of $S_1$. We retrieve the non-terminals associated within the integers $q \in T_2(u_1,u_2,v_1,v_2)$ such that the intersection of an interval $I = [a,b]$ and $[\ell,q)$ is non-empty, where $a = ([\Delta - \exp(S_1,S_2)]/\pi_1$, $b = ([\Delta - \exp(S_1,S_2) + b)/\pi_1$ and $\ell = -([|S_1| - |P_1| - \Delta_1])/\pi_1$ (See the description of the index for the definition of $\exp(S_1,S_2)$). As $\ell$ is fixed, we can implement the query via at most one binary tree search: If $b \leq \ell$, the output is empty, if $a \leq \ell \leq b$, we must output all integers, and if $\ell \leq a$, we must output all $q \geq b$. Let us now explain why the algorithm is correct. Consider a non-terminal $A$ for which $T_2(u_1,u_2,v_1,v_2)$ stores an integer $q$. By construction, $\overline{A}$ contains a relevant occurrence of $S_2$ with a split $l_2$. A position $p_1 = \lfloor \exp(A) \rfloor - l_2 - \exp(S_1,S_2) - q \cdot \pi_1$ is the leftmost occurrence of $S_1$ in $\overline{A}$ such that $p_1 \leq p_2 \leq p_1 + |S_1| - 1$ and $p_2 = \lfloor \exp(A) \rfloor - l_2 - \exp(S_1,S_2)$ the rightmost. Consequently, there is an occurrence $q_1 = \lfloor \exp(A) \rfloor - l_2 - \exp(S_1,S_2) - q \cdot \pi_1 + \Delta_1$ of $P_1$ for each $-([|S_1| - |P_1| - \Delta_1])/\pi_1 \leq q' \leq q$. The occurrence of $S_2$ implies that $q_2 = \lfloor \exp(A) \rfloor - s_2$ is an occurrence of $P_2$. We have $0 \leq q_2 - q_1 = q' \cdot \pi_1 + \exp(S_1,S_2) - \Delta \leq b$ if $\Delta - \exp(S_1,S_2) \leq q' \cdot \pi_1 \leq \Delta - \exp(S_1,S_2) + b$, which is equivalent to $[\ell,q] \cap I \neq \emptyset$. It
follows that we retrieve every non-terminal corresponding to Subcase (2.2). On the other hand, by Observation 21, the expansion of each retrieved non-terminal contains a \(b\)-close co-occurrence of \(P_1, P_2\).

**Subcase (1.6).** We argue that we have already reported all non-terminals corresponding to this subcase and there is nothing left to do. Consider a non-terminal \(A\) such that its expansion contains a relevant occurrence \(p_2\) of \(S_2\). If there are at most two occurrences \(p_1\) of \(S_1\) such that \(p_1 \leq p_2 \leq p_1 + |S_1| - 1 \leq p_2 + |S_2| - 1\), we will treat them when we query \(T_1(u_1, u_2, v_1, v_2)\) (Subcases (1.1)-(1.4)). Otherwise, by Corollary 2, \(S_1\) is periodic and there is an occurrence \(p_1'\) of \(S_1\) such that \(p_1' \leq p_2 < p_2 + \pi_1 \leq p_1 + |S_1| - 1 < p_2 + |S_2| - 1\). The non-terminals corresponding to this case are reported when we query \(T_2(u_1, u_2, v_1, v_2)\) (Subcase (2.2)).

**Time complexity.** As shown above, the algorithm reports a set \(N' \supset N\) of non-terminals and each non-terminal in \(N'\) contains a \(b\)-close co-occurrence. By Claim 16 and since the height of \(G'\) is \(h = O(\log N)\), we have \(|N'| = O(\text{occ} \log N)\). Furthermore, for a fixed pair of splits of \(P_1, P_2\), each non-terminal in \(N'\) can be reported a constant number of times. Since \(|\text{Splits}'(G', P_1)| \cdot |\text{Splits}'(G', P_2)| = O(\log^2 N)\), the total size of the output is \(|N'| \cdot O(\log^2 N) = O(\text{occ} \cdot \log^3 N)\). We therefore obtain that the running time of the algorithm is \(O(m + \log^3 N + \text{occ} \log^3 N) = O(m + (1 + \text{occ}) \log^3 N)\) as desired.

Once we have retrieved the set \(N'\), we find all \(b\)-close relevant co-occurrences for each of the non-terminals in \(N'\) using Theorem 10. In fact, our algorithm acts naively and computes all relevant co-occurrences for a non-terminal in \(N'\), and then selects those that are \(b\)-close. By case inspection, one can show that a relevant co-occurrence for a non-terminal \(A\) always consists of an occurrence of \(P_2\) that is either relevant or the leftmost in \(\text{tail}(A)\), and a preceding occurrence of \(P_1\). Intuitively, this allows to compute all relevant co-occurrences efficiently and guarantees that their number is small. Formally, we show the following claim:

**Lemma 25.** Assume that \(P_2\) is not a substring of \(P_1\). After \(O(m \log N + \log^2 N)\)-time preprocessing, the data structure of Theorem 10 allows to compute all \(b\)-close relevant co-occurrences of \(P_1, P_2\) in the expansion of a given non-terminal \(A\) in time \(O(\log^3 N \log \log N)\).

A part of the index of Christiansen et al. [8] is a pruned copy of the parse tree of \(G'\). They showed how to traverse the tree to report all occurrences of a pattern, given its relevant occurrences in the non-terminals. By using essentially the same algorithm, we can report all \(b\)-close co-occurrences in amortized constant time per co-occurrence, which concludes the proof of Theorem 1. (See Appendix C, Lemma 33.)

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**References**


A Proofs omitted from Section 2

Lemma 9. Given an RLSLP $G$ of size $g$ and height $h$. Assume that every string in a set $S$ is either a prefix or a suffix of the expansion of a non-terminal of $G$ or its reverse. The trie for $S$ can be implemented in space $O(|S|)$ to maintain the following queries in $O(m + \tau \cdot (h + \log m))$ time: Given a pattern $P$ of length $m$ and suffixes $Q_i$ of $P$, $1 \leq i \leq \tau$, find, for each $i$, the interval of strings in the (lexicographically sorted) $S$ prefixed by $Q_i$.

Proof. Let us first recall the definition of the Karp–Rabin fingerprint.

Definition 26 (Karp–Rabin fingerprint). For a prime $p$ and an $r \in \mathbb{F}_p^*$, the Karp–Rabin fingerprint $\phi$ of a string $X$ is defined as a tuple $\left( r^{|X| - 1} \mod p, r^{-|X| + 1} \mod p, \varphi_{p,r}(X) \right)$, where $\varphi_{p,r}(X) = \sum_{k=0}^{|X| - 1} S[k] r^k \mod p$.

We use the following result of Christiansen et al. [8], which builds on Belazzougui et al. [3] and Gagie et al. [16, 17].

Fact 27 ([8, Lemma 6.5]). Let $S$ be a set of strings and assume we have a data structure supporting extraction of any length-$l$ prefix of strings in $S$ in time $f_e(l)$ and computing the Karp–Rabin fingerprint $\phi$ of any length-$l$ prefix of a string in $S$ in time $f_\phi(l)$. We can then build a data structure that uses $O(|S|)$ space and supports the following queries in $O(m + f_e(m) + \tau (f_\phi(m) + \log m))$ time: Given a pattern $P$ of length $m$ and $\tau > 0$ suffixes $Q_1, \ldots, Q_\tau$ of $P$, find the intervals of strings in (the lexicographically-sorted) $S$ prefixed by $Q_1, \ldots, Q_\tau$.

It should be noted that despite using a hash function, the query algorithm is deterministic: the proof shows that $p$ and $r$ can be chosen during the construction time to ensure that there are no collisions on the substrings of the strings in $S$.

To bound $f_e$, we use [8, Lemma 6.6] which builds on Gąsieniec et al. [20] and Claude and Navarro [9].

Fact 28 ([8, Lemma 6.6]). Given an RLSLP of size $O(g)$, there exists a data structure of size $O(g)$ such that any length-$l$ prefix or suffix of $A$ can be obtained from any non-terminal $A$ in time $f_e(l) = O(l)$.
To bound \( f_h(l) \), we introduce a simple construction based on the following well-known fact:

**Fact 29.** Consider strings \( X,Y,Z \) where \( XY = Z \). Given the Karp–Rabin fingerprints of two of the three strings, one can compute the fingerprint of the third string in constant time.

**Claim 30.** Given a RLSLP \( G \) of size \( g \) and height \( h \), there exists a data structure of size \( O(g) \) that given a non-terminal \( A \) and an integer \( l \) allows to retrieve the Karp-Rabin fingerprints of the length-\( l \) prefix and suffix of \( A^r \) and \( \text{rev}(A^r) \) in time \( f_h(l) = O(h + \log l) \).

**Proof.** The claim for \( \text{rev}(A^r) \) follows for the claim for \( A^r \) by considering the grammar \( G_{rev} \), where the order of the non-terminals in each production is reversed. Below we focus on extracting the fingerprints for \( A^r \), and we further restrict our attention to prefixes of \( A^r \), the algorithm for suffixes being analogous.

The data structure consists of two sets. The first set contains the lengths of the expansions of all non-terminals in the grammar, and the second one their fingerprints.

By Fact 29 and doubling, it suffices to show an algorithm for computing the fingerprint of the length-\( l \) prefix of \( A \). Assume that \( A \) associated with a rule \( A \rightarrow BC \). If the length of \( A \) is smaller than \( l \), we return error. Otherwise, to compute the fingerprint of the length-\( l \) prefix of \( A \), we consider two cases. If \( l \leq |B| \), we recurse on \( B \) to retrieve the fingerprint of the \( l \)-length prefix of \( B \). Otherwise, we recurse on \( C \) to retrieve the fingerprint of \( C[\ldots l - |B|] \) and then compute the fingerprint of the \( l \)-length prefix of \( A \) from the fingerprints of \( B \) and \( C[\ldots l - |B|] \) in constant time by Fact 29.

For a non-terminal \( A \) associated with a rule \( A \rightarrow B^r \), we compute the fingerprint analogously. If the length of \( A \) is smaller than \( l \), we return error. Otherwise, let \( q \) be such that \( q \cdot |B| \leq l < (q + 1) \cdot |B| \). We compute the fingerprint of \( B^r \) from the fingerprint of \( B \) by applying Fact 29 \( O(1 + \log q) \) times, and the fingerprint of \( B[\ldots l - q \cdot |B|] \) recursively. We can then apply Fact 29 to compute the fingerprint of the length-\( l \) prefix of \( A \) in constant time. Note that in this case, the length of the prefix decreases by a factor at least \( q \).

If we are in a terminal \( A \), the calculation takes \( O(1) \) time (the prefix must be equal to \( A \) itself).

In total, we spend \( O(h + \log l) \) time as we recurse \( O(h) \) times, and whenever we spend more than constant time in a symbol, we charge it on the decrease in the length. The fingerprints of length-\( l \) suffixes are computed analogously.

By substituting the bounds for \( f_e(l) \) (Fact 29) and \( f_h(l) \) (Claim 28) into Fact 27, we obtain the claim of the lemma.

**B** Proofs omitted from Section 3

**Claim 31.** Given a non-terminal \( A \) of \( G' \), we can find the leftmost and the rightmost occurrences of \( P \) in \( A \) and as a corollary in \( \text{head}(A) \) and \( \text{tail}(A) \) in \( O(\log^2 N \log \log N) \) time.

**Proof.** We explain how to find the leftmost occurrence of \( P \) in \( A \); the rightmost one can be found analogously. We first check whether \( A \) contains an occurrence of \( P \) via Claim 14 in \( O(\log N \log \log N) \) time. If it does not, we can stop immediately. Below we assume that there is an occurrence of \( P \) in \( A \). Next, we check whether \( \text{head}(A) \) contains an occurrence of \( P \) via Claim 14 in \( O(\log N \log \log N) \) time. If it does, the leftmost occurrence of \( P \) in \( A \) is the leftmost occurrence of \( P \) in \( \text{head}(A) \) and we can find it by recursing on \( \text{head}(A) \). If \( \text{head}(A) \) does not contain an occurrence of \( P \), but \( A \) contains relevant occurrences of \( P \), then the leftmost occurrence of \( P \) in \( A \) is the leftmost relevant occurrence of \( P \) in \( A \) and we can
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find it in \(O(|\text{Splits}'(G', P)|) = O(\log N)\) time. Finally, if \(P\) neither occurs in \(\overline{\text{head}}(A)\) nor has relevant occurrences in \(\overline{A}\), then the leftmost occurrence of \(P\) in \(\overline{A}\) is the leftmost occurrence of \(P\) in \(\text{tail}(\overline{A})\). If \(\text{tail}(A)\) is a non-terminal \(C\), we recurse on \(C\) to find it. If \(\text{tail}(A) = B'^{-1}\) for a non-terminal \(B\), \(\text{tail}(\overline{A})\) cannot contain an occurrence of \(P\) because \(\overline{B}\) does not contain \(P\) and there are no relevant occurrences in \(A\). We recurse down at most \(h = O(\log N)\) levels, and spend \(O(\log N \log \log N)\) time per level. The claim follows.

Lemma 32. Let \(A\) be a non-terminal of \(G'\). For any position \(p\), we can find the rightmost occurrence \(q \leq p\) of \(P\) in \(\overline{A}\) and the leftmost occurrence \(q' \geq p\) of \(P\) in \(\overline{A}\) in \(O(\log^3 N \log \log N)\) time.

Proof. First we describe how to locate \(q\). Consider a node \(u\) of the parse tree of \(G'\) labeled by \(A\). The algorithm starts at \(u\) and recurses down. Let \(A'\) be the label of the current node. It computes the leftmost and rightmost occurrences in \(\overline{A}\), \(\overline{\text{head}}(A')\) and \(\overline{\text{tail}}(A')\) as well as all relevant occurrences via Claim 31. If the leftmost occurrence of \(P\) in \(\overline{A}\) is larger than \(p\), the search result is empty. Otherwise, consider two cases.

1. \(A'\) is associated with a rule \(A' \rightarrow B'P\), i.e. \(\overline{\text{head}}(A') = B', \overline{\text{tail}}(A') = C'\).
   a. If \(p \leq |\overline{B'}|\), recurse on \(B'\).
   b. Assume now that \(p > |\overline{B'}|\). If the leftmost occurrence of \(P\) in \(\overline{C'}\) is smaller than \(p\), recurse on \(C'\). Otherwise, return the rightmost relevant occurrence of \(P\) in \(\overline{A}\) if it exists else the rightmost occurrence of \(P\) in \(\overline{B'}\).
2. \(A'\) is associated with a rule \(A \rightarrow (B')^r\), i.e. \(\overline{\text{head}}(A') = B', \overline{\text{tail}}(A') = (B')^{r-1}\). Let an integer \(k\) be such that \((k - 1) \cdot |\overline{B'}| + 1 \leq p \leq k \cdot |\overline{B'}|\). The desired occurrence of \(P\) is the rightmost one of the following ones:
   a. The rightmost occurrence \(q \leq p\) of \(P\) which crosses the border between two copies of \(\overline{B'}\). To compute \(q\), we compute all relevant occurrences of \(P\) in \(\overline{A}\) and then shift each of them by the maximal possible shift \(r' \cdot |\overline{B'}|\), where \(r'\) is an integer, which guarantees that it starts before \(p\) and ends before \(|\overline{A}|\) and take the rightmost of the computed occurrences to obtain \(q\).
   b. The rightmost occurrence \(q\) of \(P\) such that for some integer \(k'\), we have \((k' - 1) \cdot |\overline{B'}| \leq q \leq q + |P| - 1 \leq k' \cdot |\overline{B'}|\) (i.e. the occurrence fully belongs to some copy of \(\overline{B'}\)). In this case, \(q\) is either the rightmost occurrence of \(P\) in the \((k - 1)\)-th copy of \(\overline{B'}\), or the rightmost occurrence of \(P\) in the \(k\)-th copy of \(\overline{B'}\) that is smaller than \(p\). In the second case, we compute \(q\) by recursing on \(B'\).

We recurse down at most \(h\) levels. On each level we spend \(O(\log^2 N \log \log N)\) time to compute the leftmost, the rightmost, and relevant occurrences and respective shifts for a constant number of non-terminals via Claim 31. Therefore, in total we spend \(O(h \cdot \log^2 N \log \log N) = O(\log^3 N \log \log N)\) time.

Locating \(q'\) is very similar and differs only in small technicalities. The algorithm starts at the node \(u\) and recurses down. Let \(A'\) be the label of the current node. We compute the leftmost and rightmost occurrences in \(\overline{A}, \overline{\text{head}}(A')\) and \(\overline{\text{tail}}(A')\) as well as all relevant occurrences via Claim 31. If the rightmost occurrence of \(P\) in \(\overline{A}\) is smaller than \(p\), the search result is empty. Otherwise, consider two cases.

1. \(A'\) is associated with a rule \(A' \rightarrow B'P\), i.e. \(\overline{\text{head}}(A') = B', \overline{\text{tail}}(A') = C'\).
   a. If \(p > |\overline{B'}|\), recurse on \(C'\).
   b. Assume now that \(p \leq |\overline{B'}|\). If the rightmost occurrence of \(P\) in \(\overline{B'}\) is larger than \(p\), recurse on \(B'\). Otherwise, return the leftmost relevant occurrence \(q\) satisfying \(q \geq p\), if it exists, and otherwise the leftmost occurrence of \(P\) in \(\overline{C'}\).
2. $A'$ is associated with a rule $A \rightarrow (B')^r$, i.e. $\text{head}(A') = B'$, $\text{tail}(A') = (B')^{r-1}$. Let an integer $k$ be such that $(k - 1) \cdot |B'| + 1 \leq p \leq k \cdot |B'|$. The desired occurrence of $P$ is the leftmost one of the following ones:

a. The leftmost occurrence $q' \geq p$ of $P$ which crosses the border between two copies of $B'$. To compute $q'$, we compute all relevant occurrences of $P$ in $\overline{A}$ and then shift each of them by the minimal possible shift $r' \cdot |B'|$, where $r'$ is an integer, which guarantees that it starts after $p$ and ends before $|\overline{A}|$ (if it exists) and take the leftmost of the computed occurrences to obtain $q$.

b. The leftmost occurrence $q'$ of $P$ such that for some integer $k'$, we have $(k' - 1) \cdot |B'| \leq q' \leq q' + |P| - 1 \leq k' \cdot |B'|$ (i.e. the occurrence fully belongs to some copy of $B'$). In this case, $q'$ is either the leftmost occurrence of $P$ in the $(k + 1)$-st copy of $B'$, or the leftmost occurrence of $P$ in the $k$-th copy of $B'$ that is larger than $p$. In the second case, we compute $q'$ by recursing on $B'$.

The time complexities are the same as for computing $q$.

\section*{C Proofs omitted from Section 4}

\begin{lemma}
Assume that $P_2$ is not a substring of $P_1$. After $O(m \log N + \log^2 N)$-time preprocessing, the data structure of Theorem 10 allows to compute all $b$-close relevant co-occurrences of $P_1, P_2$ in the expansion of a given non-terminal $A$ in time $O(\log^3 N \log \log N)$.
\end{lemma}

\textbf{Proof.} We preprocess $P_1, P_2$ in $O(m \log N + \log^2 N)$ time as explained in Theorem 10. Upon receiving a non-terminal $A$, we compute the leftmost and the rightmost occurrences of $P_1, P_2$ in $\text{head}(A)$ and $\text{tail}(A)$, as well as a set $\Pi_1$ of all relevant occurrences of $P_1$ in $\overline{A}$ and a set $\Pi_2$ of all relevant occurrences of $P_2$ in $\overline{A}$ via Claim 31. We will compute all relevant co-occurrences in $\overline{A}$, selecting those of them that are $b$-close is then trivial. As $q_1 \leq q_2$ by definition, each relevant co-occurrence $(q_1, q_2)$ of $P_1, P_2$ in $\overline{A}$ falls under one of the following categories:

1. $q_1$ is a relevant occurrence of $P_1$ in $\overline{A}$ and $q_2$ is a relevant occurrence of $P_2$ in $\overline{A}$ (i.e. $q_1 \in \Pi_1, q_2 \in \Pi_2$). To check whether a pair $q_1 \in \Pi_1, q_2 \in \Pi_2$ forms a co-occurrence of $P_1, P_2$ in $\overline{A}$, we must check whether there is an occurrence $q$ of either $P_1$ or $P_2$ between $q_1$ and $q_2$. The occurrence $q$ can only be the rightmost occurrence $r_q$ of $P_2$ in $\text{head}(A)$, the leftmost occurrence $l_q$ of $P_1$ in $\text{tail}(A)$, or an occurrence in $\Pi_1 \cup \Pi_2$. Consequently, we can find all co-occurrences in this category by merging two (sorted) sets: $\Pi_1 \cup \{l_q\}$ and $\{r_q\} \cup \Pi_2$, which can be done in $O(2 + |\Pi_1 \cup \Pi_2|)$ time.

2. $1 \leq q_1 \leq q_1 + |P_1| - 1 \leq |\text{head}(A)|$ and $|\text{head}(A)| \leq q_2 \leq q_2 + |P_2| - 1$. In this case, $q_1$ must be the rightmost occurrence of $P_1$ in $\text{head}(A)$ and $q_2$ the leftmost occurrence in $\text{tail}(A)$, $q_1 \leq q_2$, and there must be no occurrence $q \in \Pi_1 \cup \Pi_2$ such that $q_1 \leq q \leq q_2$. Therefore, if there is a co-occurrence in this category, we can retrieve it in $O(|\Pi_1 \cup \Pi_2|)$ time.

3. $q_1$ is a relevant occurrence of $P_1$ in $\overline{A}$ (i.e. $q_1 \in \Pi_1$) and $\text{head}(A) \leq q_2 \leq q_2 + |P_2| - 1$. In this case, $q_1$ must be the rightmost occurrence in $\Pi_1$ and $q_2$ the leftmost occurrence of $P_2$ in $\text{tail}(A)$, and there should be no occurrence from $\Pi_2$ between $q_1$ and $q_2$. Therefore, if there is a co-occurrence in this category, we can find it in $O(|\Pi_1 \cup \Pi_2|)$ time.

4. $q_1 \leq q_1 + |P_1| - 1 \leq |\text{head}(A)|$ and $q_2$ is a relevant occurrence of $P_2$ in $\overline{A}$ (i.e. $q_2 \in \Pi_2$). First, consider the leftmost occurrence in $\overline{A}$. We find the rightmost occurrence $q_1 \leq q_2$ of $P_1$ in $\overline{A}$ via a predecessor query. The pair $(q_1, q_2)$ is a co-occurrence if the rightmost occurrence of $P_2$ in $\text{head}(A)$ is smaller than $q_1$, which can be checked in constant time. Second, we consider the remaining occurrences in $\Pi_2$. Let $q'_2$ be the leftmost one.
We begin by computing the preceding occurrence $q'_1$ of $P_1$ via a predecessor query and if $q_2 ≤ q'_1$, output the resulting co-occurrence. If $Π_2 = {q_2, q'_2}$, we are done. Otherwise, by Corollary 2, the occurrences in $Π_2 \setminus \{q_2\}$ form an arithmetic progression with difference equal to the period of $P_2$ (as all of them contain the position $\text{head}(A)$). Furthermore, as $P_1$ does not contain $P_2$, the occurrence of $P_1$ preceding $q'_2$ belongs to the periodic region formed by the relevant occurrences of $P_2$. Therefore, all the remaining co-occurrences can be obtained from the co-occurrence for $q'_2$ by shifting them by the period. In total, this step takes $O(|Π_2| + \log^5 N \log \log N)$ time.

Lemma 33. Assume that $P_2$ is not a substring of $P_1$. One can compute all $b$-close co-occurrences of $P_1, P_2$ in $S$ in time $O(m + (1 + \text{occ}) \cdot \log^4 N \log \log N)$.

Proof. During the preprocessing, we prune the parse tree: First, for each non-terminal $B$, all but the first node labeled by $B$ in the preorder is converted into a leaf and its subtree is pruned. For each node $v$ labeled by a non-terminal $B$, we store $\text{anc}(v)$, the nearest ancestor $u$ of $v$ labeled by $A$ such that $u$ is the root or $A$ labels more than one node in the pruned tree. Second, for every node labeled by a non-terminal $A$ associated with a rule $A \rightarrow B^k$, we replace its $k - 1$ rightmost children with a leaf labeled by $B^{k-1}$. We call the resulting tree the pruned parse tree and for each node $v$ labeled by a non-terminal $B$ store $\text{next}(v)$, the next node labeled by $B$ in preorder, if there is one. As every non-terminal labels at most one internal node of the pruned parse tree and every node has at most two children, it occupies $O(g')$ space.

When the algorithm of Lemma 24 outputs $A \in N'$, we compute all relevant co-occurrences $(q_1, q_2)$ in $A$ in time $O(\log^5 N \log \log N)$ using Lemma 25 and select those which satisfy $q_2 - q_1 ≤ b$.

Fix a $b$-close relevant co-occurrence $(q_1, q_2)$ in $A$. If $A$ is associated with a rule $A \rightarrow BC$, construct a set $\text{occ}(A) := \{(q_1, q_2)\}$, and otherwise if $A$ is associated with a rule $A \rightarrow B^k$,

$$\text{occ}(A) := \{(q_1 + i \cdot |\overline{B}|, q_2 + i \cdot |\overline{B}|) : 0 ≤ i ≤ \left|\left(\left|\overline{A}\right| - q_2 - |P_2| + 1\right)/|\overline{B}|\right|\}$$

Suppose that $A$ labels nodes $v_1, v_2, \ldots, v_k$ of the unpruned parse tree of $G'$ (by construction $v_1$ is not pruned and we assimilate it to the corresponding node in the pruned parse tree). If $W$ is a set of co-occurrences, denote for brevity $W + \delta = \{(q_1 + \delta, q_2 + \delta) : (q_1, q_2) \in W\}$. Below we show an algorithm that generates a set $S = \cup_i \text{occ}(A) + \text{off}(v_i)$ that contains all secondary $b$-close co-occurrences due to $(q_1, q_2)$.

We traverse the pruned parse tree, while maintaining a priority queue. The queue is initialized to contain the first node in the preorder labeled by $A$ together with $\text{occ}(A)$. Until the priority queue is empty, pop a node $v$ and a set $W$ of co-occurrences of $P_1, P_2$ in the expansion of its label, and perform the following steps:

- **Reporting step**: If $v$ is the root, report $W$;
- **Next node step**: If $\text{next}(v)$ is defined, push $(\text{next}(v), W + \text{off}(\text{next}(v)) - \text{off}(v))$;
- **Sibling step**: If $v$ is labeled by a non-terminal $B$ and its sibling by $B^k$, for some integer $k$, then $W := \cup_{0 \leq i ≤ k} W + i \cdot |\overline{B}|$
- **Ancestor step**: Push to the queue $(\text{anc}(v), W + \text{off}(\text{anc}(v)) - \text{off}(v))$.

By construction and as every node is connected with the root by a path of anc links, the algorithm generates each co-occurrence in $S$ exactly once. The time complexity follows: The algorithm of Lemma 24 takes $O(m + (1 + \text{occ}) \cdot \log^5 N)$ time; applying Lemma 25 to every non-terminal in $N'$ takes $O(\text{occ} \cdot \log^4 N \log \log N)$ time; and maintaining the queue and reporting the co-occurrences takes $O(\text{occ})$ time as at every step we can charge the time needed to update the queue on newly created co-occurrences.
Order-Preserving Squares in Strings

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Abstract

An order-preserving square in a string is a fragment of the form $uv$ where $u \neq v$ and $u$ is order-isomorphic to $v$. We show that a string $w$ of length $n$ over an alphabet of size $\sigma$ contains $O(\sigma n)$ order-preserving squares that are distinct as words. This improves the upper bound of $O(\sigma^2 n)$ by Kociumaka, Radoszewski, Rytter, and Waleń [TCS 2016]. Further, for every $\sigma$ and $n$ we exhibit a string with $\Omega(\sigma n)$ order-preserving squares that are distinct as words, thus establishing that our upper bound is asymptotically tight. Finally, we design an $O(\sigma n)$ time algorithm that outputs all order-preserving squares that occur in a given string and are distinct as words. By our lower bound, this is optimal in the worst case.

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1 Introduction

A natural definition of repetitions in strings is that of squares, which are fragments of the form $uu$, where $u$ is a string. The study of repetitions in strings goes back at least to the work of Thue from 1906 [28], who constructed an infinite square-free word over the ternary alphabet. Since then, multiple definitions of repetitions have been proposed and studied, with the basic question being focused on analyzing how many such repetitions a string of length $n$ can contain. Of course, any even-length fragment of the string $a^n$ is a square, therefore we would like to count distinct squares. Using a combinatorial result of Crochemore and Rytter [5], Fraenkel and Simpson [10] proved that a string of length $n$ contains at most $2n$ distinct squares (also see a simpler proof by Ilie [17]). They also provided an infinite family of strings of length $n$ with $n - o(n)$ distinct squares. For many years, it was conjectured that the right upper bound is actually $n$. Interestingly, a proof of the conjecture for the binary alphabet would imply it for any alphabet [24]. Very recently, after a series of improvements on the upper bound [7,18,23,27], the conjecture has been finally resolved by Brlek and Li [1], who showed an upper bound of $n - \sigma + 1$, where $\sigma$ is the size of the alphabet.

For many of the applications, it seems more appropriate to work with different definitions of equality, giving us different notions of squares. Three interesting examples are (1) Abelian squares [6,8,9,16,19–21,26] (also called Jumbled squares) are of interest in natural language
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processing applications and in other domains where the classifications strongly depend on feature sets distribution, as opposed to feature sequences distributions. (2) Parameterized squares [20] are considered in applications for finding identical sections of code. (3) Order-preserving squares [4,13,20] could be used in applications of stock price analysis and musical melody matching.

The combinatorial properties of the three types of squares were studied by Kociumaka et al. [20]. Given a string of length $n$ over an alphabet of size $\sigma$, first the authors bounded the number of abelian squares that are distinct as words by $\Theta(n^2)$. Second, bounded the number of parameterized squares that are distinct as words by $O((\sigma!)^2 n)$ and bounded the number of nonequivalent parameterized squares (see definition within) by $O(\sigma n)$. Third, the authors provided $O(\sigma^2 n)$ bound for the number of order-preserving squares that are distinct as words.

From an algorithmic perspective, various algorithms were proposed for computing abelian squares and order-preserving squares in a string of length $n$. Cummings and Smyth [6] proposed an $\Theta(n^2)$ time algorithm for computing all substrings that consist of a concatenation of two or more abelian-equivalent substrings. Kociumaka et al. [21] proposed an algorithm for computing the longest, the shortest, and the number of all abelian squares in $O(n^2/\log^2 n)$ time using linear space. Gourdel et al. [13] proved that all nonshiftable order-preserving squares (see definition within) can be computed in $O(n \log n)$ time. Additionally, Crochemore et al. [4] proposed the incomplete order-preserving suffix tree (see details within), denoted by $T$, that enables order-preserving pattern matching queries in time proportional to the pattern length. The suffix tree $T$ can be constructed in $O(n \log \log n)$ expected time and $O(n \log^2 \log n/\log \log \log n)$ worst-case time. Moreover, the authors proved that using $T$, all occurrences of order-preserving squares can be computed in $O(n \log n + occ)$ time, where $occ$ is the total number of occurrences of order-preserving squares. Note that, the number of all occurrences of order-preserving squares might be unreasonably high. In particular, every regular square is considered to be an order-preserving square, hence $a^n$ contains $\Theta(n^2)$ occurrences of order-preserving squares. Henceforth, a more natural approach is to generate only order-preserving squares that are distinct as words.

**Our results.** In this paper, we focus on order-preserving squares. Same-length strings $u$ and $v$ over an ordered alphabet are order-isomorphic, denoted $u \approx v$, when the order between the characters at the corresponding positions is the same in $u$ and $v$. For example, the strings $u = acb$ and $v = azd$ are order-isomorphic, assuming $a < b < c < d < z$. In this paper, order-preserving squares are strings of the form $uv$, where $u \approx v$ and additionally $u \neq v$.

The main result of our paper is that the number of order-preserving squares in a string of length $n$ over an alphabet of size $\sigma$ is $O(\sigma n)$. This improves the bound of $O(\sigma^2 n)$ by Kociumaka et al. [20]. We stress that in our definition of an order-preserving square, we require that $u \neq v$, while Kociumaka et al. [20] counted fragments of the form $uv$, where $u \approx v$, that are distinct as words. We believe that our definition is more natural in the context of this paper. At the same time, by the result of Brlek and Li [1] a string of length $n$ contains less than $n$ fragments of the form $uv$ that are distinct as words, thus our result implies that the number of fragments $uv$ such that $u \approx v$ that are distinct as words is also $O(\sigma n)$. We complement our upper bound by designing, for each $\sigma$, an infinite family of strings of length $n$ over an alphabet of size $\sigma$ containing $\Theta(\sigma n)$ such fragments. We begin with describing the lower bound in Section 3, and then present the upper bound in Section 4.

**Theorem 1.** The number of order-preserving squares in a string of length $n$ over an alphabet of size $\sigma$ is $O(\sigma n)$, and this bound is asymptotically tight even if we only consider order-preserving squares that are distinct as words.
Next, we design an algorithm for reporting all order-preserving squares in a given string of length \( n \) over an alphabet of size \( \sigma \) in \( O(\sigma n) \) time, which (by our lower bound) is asymptotically optimal in the worst case. We again stress that in our definition of an order-preserving square, we require that \( u \neq v \). However, all fragments of the form \( uw \) that are distinct as words can be reported in \( O(\sigma n) \) time using the algorithm of Gusfield and Stoye \([14]\)\(^1\). Thus, for \( \sigma = o(\log n) \), this resolves one of the open questions by Crochemore et al. \([4]\), who asked if there is an \( o(n \log n) \) time algorithm for finding the longest order-preserving square. This is described in Section 5.

**Theorem 2.** All order-preserving squares in a string of length \( n \) over an alphabet of size \( \sigma \) can be found in \( O(\sigma n) \) time.

**High-level description of our techniques.** For the lower bound, first, we consider the increasing string \( w = 123\ldots n \) where \( \sigma = n \). Clearly, any even-length fragment is an order-preserving square thus producing the maximum number, i.e. \( \Omega(n^2) = \Omega(\sigma n) \), of order-preserving squares in a string of length \( n \). To decrease the size of the alphabet \( \sigma \), we replace \( w \) with a non-decreasing string \( w = 11\ldots 2\ldots \sigma \ldots \sigma \), where each character is repeated the same number of times. We exhibit \( \Omega(\sigma n) \) order-preserving squares in \( w \) that are distinct as words. See Section 3 for more details.

For the upper bound, we build on the insight by Kociumaka et al. \([20]\), where the high-level strategy is to consider each suffix of \( w \) separately. For each suffix and an alphabet character, they considered the leftmost occurrence of this character within the suffix. Thus, there are at most \( \sigma \) leftmost occurrences in each suffix. For a fixed suffix, they considered all of its prefixes as possible order-preserving squares \( uv \). Next, they showed that, because \( u \neq v \), the order-preserving square \( uv \) is defined by a pair (or pairs) of leftmost occurrences such that one occurrence belongs to \( u \), and the other one belongs to \( v \) at the same relative position, where the length of \( uv \) is twice the difference between the leftmost occurrences. For example, let \( acbadxyz \) be the suffix, then the pair of positions 2 and 5 are leftmost occurrences defining the order-preserving square \( acbadx \) of length 6 that is a prefix of the given suffix. Note that, also 3 and 6 are leftmost occurrences defining the same order-preserving square \( acbadx \).

Thus, as a result, they upper bounded the number of order-preserving squares being a prefix of the considered suffix by \( (\frac{\sigma}{2})n \), so \( (\frac{\sigma}{2})n \) in total.

In this paper, we adopt a similar approach, by separately upper bound the number of order-preserving squares that are prefixes of a suffix of the input string \( w \). However, our goal is to show that there are only \( O(\sigma) \) such prefixes, so \( O(\sigma n) \) in total. To this end, we first partition the order-preserving squares into groups. Let \( O_k \) the set of all order-preserving squares \( uv \) such that \( 2^k \leq |uv| < 2^{k+1} \). Similarly, we partition the leftmost occurrences into groups. Let \( L_k \) the set of all leftmost occurrences \( i \) such that \( 2^k \leq i < 2^{k+1} \). Now, our strategy is to show that if \( |O_k| \) is larger than some fixed constant then \( |O_k| = O(|L_k|) \). The structure of the argument is as follows. We first observe that two order-preserving squares \( uv \) and \( u'v' \) imply that \( |u| - \Delta, \) where \( \Delta = |u'| - |u| \), is a so-called order-preserving border of \( u \). We write \( u = b_1b_2\ldots b_f b_{f+1}, \) where \( |b_1| = |b_2| = \ldots |b_f| = \Delta \) and \( |b_{f+1}| < \Delta, \) and by carefully choosing \( uv \) and \( u'v' \) from \( O_k \) conclude that \( b_2 \) contains a leftmost occurrence and \( f \) is proportional to \( |O_k| \). Then, we argue that \( b_2 \) containing a leftmost occurrence implies

---

\(^1\) They only claim \( O(n) \) time for fixed alphabets, however a closer look at the algorithm reveals that there are 3 phases: the first phase takes \( O(n) \) time (Theorem 6 and Lemma 7), the second phase also takes \( O(n) \) time (Section 4), and the third phase takes \( O(\sigma n) \) time (Lemma 11). Additionally, the algorithm assumes that the suffix tree is constructed in \( O(n) \) time, for larger alphabets this increases to \( O(\sigma n) \).
that, in fact, every $b_j$ contains a leftmost occurrence, and thus $|O_k| = \mathcal{O}(|L_k|)$. Summing this over all $k$, and separately considering all $k$ such that $|O_k|$ is less than the fixed constant, we are able to conclude that $\sum_k |O_k| = \sum_k \mathcal{O}(|L_k|) < \mathcal{O}(\sigma)$. See Section 4 for more details.

To obtain an efficient algorithm for reporting all order-preserving squares, we apply the order-preserving suffix tree as defined by Crochemore et al. [4]. This structure allows us to check if $w[i..i+2\ell-1]$ is an order-preserving square by checking if the LCA of two leaves is at string depth at least $\ell$. First, we need to show how to construct the order-preserving tree in $\mathcal{O}(\sigma n)$ time. Second, we extend the above reasoning to efficiently generate only $\mathcal{O}(\sigma n)$ fragments that are then tested for being an order-preserving square in constant time each. While the underlying argument is essentially the same as when bounding the number of order-preserving squares, it needs to be executed differently for the purpose of an efficient implementation. See Section 5 for more details.

2 Preliminaries

Let $\Sigma = \{1, \ldots, \sigma\}$ be a fixed finite alphabet of size $\sigma$. Let $|s|$ denote the length of a string $s$. For a string $s$, the character at position $i$ of $s$ is denoted by $s[i]$, and $s[i..j]$ is the fragment of $s$ starting at position $i$ and ending at position $j$. We call two strings $u$ and $v$ order-isomorphic, denoted by $u \approx v$, when $|u| = |v|$ and, for each $i, j$, we have $u[i] \leq u[j]$ if and only if $v[i] \leq v[j]$. The concatenation of two strings $u$ and $v$ is denoted by $uv$. A string of the form $uv$ is called an order-preserving square, or op-square, when $u \neq v$ and $u \approx v$. We call $u$ its left arm and $v$ its right arm. We stress that a regular square, that is, a string of the form $xx$, is not an op-square. Two op-squares $uv$ and $u'v'$ are distinct as words if and only if $uv \neq u'v'$.

A trie is a rooted tree, with every edge labeled with a single character and edges outgoing from the same node having distinct labels. A node $u$ of a trie represents the string obtained by reading the labels on the path from the root to $u$. A compacted trie is obtained from a trie by replacing maximal paths consisting of nodes with exactly one child with single edges labeled by the concatenation of the labels of the edges on the path. A suffix tree $T$ of a string $w$ is a compacted trie whose leaves correspond to the suffixes of $w$. The string depth of a node $u$ of $T$ is the length of the string that it corresponds to. An explicit node of $T$ is simply a node of $T$. An implicit node of $T$ is a node of the non-compacted trie corresponding to $T$, or in other words a location on an edge of $T$.

Next, we need some definitions specific to order-isomorphism. Following Kubica et al. [22], we call $b$ an op-border of a string $s[1..n]$ when $s[1..b] \approx s[n-b+1..n]$. Following Gourdel et al. [13] (and Matsuoka et al. [25]), we call $p$ an (initial) op-period of $s[1..n]$ when $s = b_1b_2\ldots b_f b_{f+1}$ with $|b_1| = |b_2| = \ldots = |b_f| = p$ and $|b_{f+1}| < p$ (so $f = \lfloor n/p \rfloor$). $b_1 \approx b_2 \approx \ldots \approx b_f$ and $b_1[1..|b_{f+1}|] \approx b_2[1..|b_{f+1}|] \approx \ldots \approx b_f[1..|b_{f+1}|] \approx b_{f+1}$. $b_1, b_2, \ldots, b_f$ are called the blocks defined by $p$ in $s$, while $b_{f+1}$ (possibly empty) is called the incomplete block. While in the classical setting $p$ is a period of $s[1..n]$ if and only if $n-p$ is a border of $s[1..n]$, in the order-preserving setting, we only have an implication in one direction. For example, the string $a\text{f}ic\text{d}g\text{d}b\text{e}h$ has an op-period 3 while 6 is not its op-border.

\begin{proposition}
If $b$ is an op-border of $s[1..n]$ then $n-b$ is an initial op-period of $s[1..n]$.
\end{proposition}

\begin{proof}
Let $p = n-b$ and $f = \lfloor n/p \rfloor$. We represent $s[1..n]$ as $s = b_1b_2\ldots b_f b_{f+1}$ with $|b_1| = |b_2| = \ldots = |b_f| = p$ and $|b_{f+1}| < p$. By $b$ being an op-border of $s[1..n]$, we have $s[1..b] \approx s[n-b+1..n]$, so $s[1..n-p] \approx s[p+1..n]$. We observe that $s[1..n-p] = b_1b_2\ldots b_{f-1}b_f[1..|b_{f+1}|]$ and $s[p+1..n] = b_2b_3\ldots b_f b_{f+1}$. Then, $b_1b_2\ldots b_{f-1}b_f[1..|b_{f+1}|] \approx
For every $i = 1, 2, \ldots, f$, we obtain $b_1 \approx b_2 \approx b_3 \approx \ldots \approx b_{f-1} \approx b_f$ and $b_1 \approx b_2 \approx b_3 \approx \ldots \approx b_{f-1} \approx b_f$, so $p = n - b$ is indeed an initial op-period of $s[1..n]$.

Due to Proposition 3, if $b$ is an op-border of $s[1..n]$ then $s[1..n] = b_1 b_2 \ldots b_f b_{f+1}$, where $b_1 b_2 \ldots b_f | b_{f+1}| \approx b_2 b_3 \ldots b_f b_{f+1}$, $|b_1| = |b_2| = \ldots = |b_f| = n - b$ and $|b_{f+1}| < p$ (so $f = [n/(n - b)]$, $b_1 \approx b_2 \approx \ldots \approx b_f$ and $b_1 \approx b_2 \approx b_3 \approx \ldots \approx b_{f-1} \approx b_f$, so $p = n - b$ is indeed an initial op-period of $s[1..n]$.

We will say that these blocks are defined by $b$.

### Lower Bound

Recall that $\Sigma = \{1, \ldots, \sigma\}$. We define a string $w = 11 \ldots 1222 \ldots 2\sigma\ldots\sigma$, that is, a concatenation of $\sigma$ blocks, each consisting of $k$ repetitions of the same character. We note that $|w| = \sigma k$. For $i = 1, 2, \ldots, [\sigma/2]$, we consider all fragments of $w$ of length $2ik$ starting at positions $j = 1, 2, \ldots, |w| - 2ik + 1$. For $j = 1$ mod $k$, the fragment is a concatenation of $2i$ blocks, each block consisting of $k$ repetitions of the same character. For $j \neq 1$ mod $k$, the fragment starts with $r \in [1, k - 1]$ repetitions of the same character, then $2i - 1$ blocks, each block consisting of $k$ repetitions of the same character, and finally $\ell = k - r$ repetitions of the same character. See Figure 1.

![Figure 1](image)

Figure 1 The red box corresponds to an op-square of length 10 containing 3 different characters. The blue box corresponds to an op-square of length 20 containing 5 different characters.

Each such fragment is an op-square. For $j = 1$ mod $k$, both the left and the right arm consist of $i$ blocks consisting of $k$ repetitions of character $c, c + 1, \ldots, c + i - 1$. For $j \neq 1$ mod $k$, both the left and the right arm consist of first $r$ repetitions of character $c$, then $i - 1$ blocks consisting of $k$ repetitions of characters $c + 1, c + 2, \ldots, c + i - 2$, and then finally $\ell$ repetitions of character $c + i - 1$. Thus, the left and the right arm are always order-isomorphic. Further, for every choice of $i$ and the starting position we obtain a different word, as two such fragments of the same length either start with different characters or differ in the length of the first block of the same character.

Now, we analyze the number of such op-squares in $w$. By considering every $1 \leq i \leq [\sigma/2]$ and starting position $1, 2, \ldots, |w| - 2ik + 1$, we obtain that the number of op-squares in $w$ is at least:

\[
\sum_{i=1}^{[\sigma/2]} (|w| - 2ik + 1) = \sum_{i=1}^{[\sigma/2]} (\sigma k - 2ik + 1) = \frac{\sigma}{2} \cdot (\sigma k - k([\sigma/2] + 1) + 1) \\
\geq \frac{\sigma}{2} \cdot (k([\sigma/2] - 1) + 1).
\]

For $\sigma \geq 3$, this is at least $\sigma^2 k/12 = \sigma n/12$ for any $k$. For $\sigma = 1, 2$, we additionally assume $k \geq 2$ and count op-squares of the form $1^2$, there are $[k/2] \geq n/6 \geq \sigma n/12$ of them. Thus, in either case for every $k \geq 2$ we obtain a string of length $n = k\sigma$ over $\Sigma$ containing $\sigma n/12$ op-squares that are distinct as words.

*Theorem 4.* For any alphabet $\Sigma = \{1, 2, \ldots, \sigma\}$, there exists an infinite family of strings of length $n = k\sigma$ over $\Sigma$ containing $\Omega(\sigma n)$ op-squares distinct as words.
4 Upper bound

Our goal in this section is to upper bound the number of op-squares in a given string \( w \) of length \( n \) over the alphabet \( \Sigma = \{1, \ldots, \sigma\} \). Recall that \( uv \) is an op-square when \( u \neq v \) and \( u \approx v \). We will show that this number is \( O(\sigma n) \). As explained in the introduction, by the result of Brlek and Li [1], the number of regular squares, that is, fragments of the form \( uu \) that are distinct as words, is less than \( n \). Thus, our result in fact allows us to upper bound the number of fragments of the form \( uv \), where \( u \approx v \), that are distinct as words by \( O(\sigma n) \).

We consider each suffix of \( w \) separately. For each suffix \( w[i..n] \), we will upper bound the number of prefixes of \( w[i..n] \) that are op-squares by \( O(\sigma) \). Therefore, to avoid cumbersome notation in the remaining part of this section we will assume that we have a string \( s \) of length \( m \) over the alphabet \( \Sigma = \{1, \ldots, \sigma\} \), and we want to upper bound the number of op-squares \( uv \) that are prefixes of \( s \) by \( O(\sigma) \). See Figure 2.

\[
S = a_1 a_2 a_3 \cdots \cdots \cdots \cdots \cdots a_m
\]

![Figure 2](image)

**Figure 2** Green prefixes of \( s \) are op-squares.

Kociumaka et al. [20] observed that every op-square \( uv \) that is a prefix of \( s \) can be obtained as follows (recall that in our definition \( u \neq v \)). We call position \( i \) a lefmost occurrence and \( s[i] \) a leftmost character when \( s[j] \neq s[i] \) for every \( j < i \). Then, there exists \( i \) and \( j \) such that both \( i \) and \( j \) are lefmost occurrences, where \( i \) belongs to \( u \) and \( j \) belongs to \( v \), and further \( |u| = j - i \). More formally:

**Proposition 5** ([20, Lemma 4.2 and Corollary 4.3]). We can construct an injective function \( g \) mapping op-squares that are prefixes of \( s \) to 2-element subsets of the alphabet as follows. We choose the smallest \( i \) belonging to \( v \) such that \( s[i] = a \) does not occur in \( u \), and let \( s[i - |u|] = b \) be its counterpart in \( u \), then set \( g(uv) = \{a, b\} \). Both \( i \) and \( i - |u| \) are lefmost occurrences.

We split all op-squares that are prefixes of \( s \) into groups. Let \( O_k \) denote the group of op-squares that are prefixes of \( s \) having length at least \( 2^k \) and at most \( 2^{k+1} - 1 \):

**Definition 6.** \( O_k = \{uv | u \neq v \text{ and } u \approx v \text{ and } 2^k \leq |uv| < 2^{k+1}\} \) for \( 0 \leq k \leq \log m \).

In other words, we split \( s \) into consecutive ranges of exponentially increasing lengths, such that the \( k \)-th range is of length \( 2^k - 1 \), starts at position \( 2^k \) and ends at position \( 2^{k+1} - 1 \) in \( s \) (where \( 0 \leq k \leq \log m \) and the final range may not be complete when \( m < 2^{k+1} - 1 \)). Then, the set \( O_k \) consists of op-squares that end in the \( k \)-th range. See Figure 3.

The number of op-squares \( uv \) that are prefixes of \( s \) is \( \sum_{k=0}^{\log m} |O_k| \). In order to upper bound the sum, we will separately upper bound the size of each group. We first need some propositions.

**Proposition 7.** For any \( \{uv, u'v'\} \in O_k \) such that \( |u| < |u'| \) and \( \Delta = |u'| - |u|, |u| - \Delta\) is an op-border of both \( u \) and \( v \).
Because $u \approx v$ it is enough to show that $|u| - \Delta$ is an op-border of $u$. By the assumption that both $uv$ and $u'v'$ are op-squares we have:

$$u[1..|u| - \Delta] = u'[1..|u| - \Delta] \approx v'[1..|u| - \Delta] = v[\Delta + 1..|u|] \approx u[\Delta + 1..|u|].$$

See Figure 4.

In the remaining part of this section, we will often consider $\{uv, u'v'\} \in O_k$ such that $|u| < |u'|$ and $\Delta = |u'| - |u|$. Then, by Proposition 7, we know that $|u| - \Delta$ is an op-border of $u$, and thus by Proposition 3 $\Delta$ is an initial op-period of $u$. Hence, $u$ can be represented as a concatenation of $f = \lfloor |u|/\Delta \rfloor$ blocks $b_1, b_2, \ldots, b_f$ and one incomplete block $b_{f+1}$, where $|b_1| = |b_2| = \ldots = |b_f| = \Delta$ and $|b_{f+1}| < \Delta$, such that $b_1 b_2 \ldots b_f [1..|b_{f+1}|] \approx b_2 b_3 b_f b_{f+1}$, $b_1 \approx b_2 \approx \ldots \approx b_f$ and $b_1 [1..|b_{f+1}|] \approx b_2 [1..|b_{f+1}|] \approx \ldots \approx b_f [1..|b_{f+1}|] \approx b_{f+1}$. See Figure 5.

For brevity, in the remaining part of the paper we will describe this situation by saying that $\{uv, u'v'\} \in O_k$ define blocks $b_1, b_2, \ldots, b_f, b_{f+1}$.

\begin{itemize}
\item Figure 3 Green prefixes of $s$ are op-squares ending in the $k$-th range. The red line illustrates the ranges.
\item Figure 4 The green lines correspond to $uv$ and $u'v'$. The blue line corresponds to $u[1..|u| - \Delta]$. The orange line corresponds to $u[\Delta + 1..|u|]$. The purple line corresponds to $v[\Delta + 1..|u|]$.
\item Figure 5 Blocks defined by $\{uv, u'v'\}$ in $u$.
\end{itemize}

\textbf{Proposition 8.} If $|O_k| \geq 3$ then there exist $\{uv, u'v', u''v''\} \in O_k$ such that $0 < |u''| - |u|, |u''| - |u'| < 2^k/|O_k| - 2$.

\textbf{Proof.} The length of every op-square in $O_k$ belongs to $[2^k, 2^{k+1})$, thus the length of its left arm falls within $[2^{k-1}, 2^k)$. Let $O_k = \{u_{1v_1}, u_{2v_2}, \ldots, u_{tv_t}\}$ with $|u_1| < |u_2| < \ldots < |u_t|$. Then, for some $i \in \{1, 2, \ldots, \lceil (t-1)/2 \rceil\}$ we must have $|u_{2i+1}| < |u_{2i-1}| + 2^{k-1}/\lceil (t-1)/2 \rceil$ (as otherwise we would have $u_i \geq u_1 + 2^{k-1}$). The sought op-squares are $u_{2i-1}u_{2i-1}, u_{2i}u_{2i}, u_{2i+1}u_{2i+1}$ because:
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\[ |u_{21}| - |u_{21-1}|, |u_{21+1}| - |u_{21}| < |u_{21+1}| - |u_{21-1}| < 2^{k-1}/(\ell - 1/2) \]
\[ \leq 2^{k-1}/(\ell/2 - 1) = 2^k/(|O_k| - 2). \]

With all the propositions in hand, we are now ready for the technical lemmas. Our goal is to upper bound \( \sum_k |O_k| \) by the number of leftmost occurrences. To this end, we need to show that, if some \( O_k \) is large then there are many leftmost occurrences in some range. This will be done by applying the following reasoning to the three op-squares chosen by applying Proposition 8. In the following, whenever we refer to a leftmost occurrence in block \( b \) we mean a leftmost occurrence falling within the positions in block \( b \).

\[ \textbf{Lemma 9.} \text{If } |O_k| \geq 3 \text{ then for any } \{uv, u'v', u''v''\} \in O_k \text{ where } |u| < |u'| < |u''| \text{ such that } \{uv, u'v'\} \text{ defines } b_1, \ldots, b_f, b_{f+1} \text{ and } \{u''v', u''v''\} \text{ defines } b'_1, \ldots, b'_f, b'_{f+1} \text{ there is a leftmost occurrence in block } b_j \text{ such that } j \neq 1 \text{ or there is a leftmost occurrence in block } b'_j, \text{ such that } j' \neq 1. \]

\[ \textbf{Proof.} \text{Let } \Delta = |u'| - |u| \text{ be the length of every block } b_j \text{ and } \Delta' = |u''| - |u'| \text{ be the length of every block } b'_f. \text{ By Proposition 5, we know that there must be a leftmost occurrence } i \text{ that falls within } u' \text{ and its corresponding leftmost occurrence } i + |u'| \text{ that falls within } v'. \text{ Thus, we have found a leftmost occurrence } i + |u'| \text{ falls within } v. \text{ To verify this, we calculate:} \]
\[ i + |u'| \leq \Delta' + |u'| = |u''| - |u'| + |u'| = |u''| < 2^k \leq |uv|. \]

We have established that \( i + |u'| \) is a leftmost occurrence and falls within \( v \). Thus, \( s[i'] \neq s[i + |u'|] \text{ for every } i' \in [i, i + |u'|). \text{ Because } u \approx v, \text{ this then implies that } s[i' - |u|] \neq s[i + |u'| - |u|] \text{ for every } i' \in [|u| + 1, i + |u'|). \text{ Thus, } i + |u'| - |u| \text{ is also a leftmost occurrence. We claim that } s[i + |u'| - |u|] \text{ cannot belong to } b_1. \text{ To verify this, we calculate:} \]
\[ i + |u'| - |u| \geq 1 + |u'| - |u| = 1 + \Delta. \]

Thus, we have found a leftmost occurrence \( i + |u'| - |u| \) that falls within \( u \) and belongs to a block \( b_j \) with \( j \neq 1 \). See Figure 6.

\[ \textbf{Figure 6} \text{ The red points correspond to the leftmost occurrences considered in the proof of Lemma 9.} \]

Next, we show that if \( \{uv, u'v'\} \in O_k \text{ define blocks } b_1, b_2, \ldots, b_f, b_{f+1} \text{ such that there is a leftmost occurrence in block } b_j \text{ for some } j \neq 1 \text{ then, in fact, there is a leftmost occurrence in every block } b_j. \text{ This reasoning is done in two steps.} \]

\[ \textbf{Lemma 10.} \text{Let } b \text{ be an op-border of } u = s[1..|u|] \text{ that defines blocks } b_1, b_2, \ldots, b_f, b_{f+1}, \text{ and assume that there is a leftmost occurrence in block } b_j, \text{ for some } j \in [1, f + 1]. \text{ Then there is a leftmost occurrence in every block } b_1, b_2, \ldots, b_j. \]
Proof. Let $\Delta = |b_1| = |b_2| = \ldots = |b_f|$ and $|b_{f+1}| < \Delta$. By induction, it is enough to show that if there is a leftmost occurrence in block $b_j$ for some $j \geq 2$ then there is a leftmost occurrence in block $b_{j-1}$. Let $i$ be a leftmost occurrence that belongs to $b_j$. Then $u[i'] \neq u[i]$ for every $i' \in [1,i)$. Because $b_1b_2\ldots b_{j-1}b_j[1..|b_{f+1}|] \approx b_2b_3\ldots b_fb_{f+1}$, this implies $u[i'-\Delta] \neq u[i-\Delta]$ for every $i' \in [\Delta+1, i)$. But then $i - \Delta$ is also a leftmost occurrence, and it belongs to $b_{j-1}$ as required. See Figure 7.

![Figure 7](image)

Each red point corresponds to a leftmost character at the same relative position in every block $b_1, b_2, \ldots, b_f$.

Lemma 11. Let $b$ be an op-border of $u = s[1..|u|]$ that defines blocks $b_1, b_2, \ldots, b_f, b_{f+1}$, and assume that there is a leftmost character in block $b_2$. Then there is a leftmost occurrence in every block $b_1, b_2, \ldots, b_f$.

Proof. Let $\Delta = |b_1| = |b_2| = \ldots = |b_f|$ and $|b_{f+1}| < \Delta$. By assumption, there is a leftmost character $s[i]$ in block $b_2$, where $i \in [\Delta + 1, 2\Delta]$. Our goal is to show that there is a leftmost occurrence in every block $b_1, b_2, \ldots, b_f$.

Because $b_1b_2\ldots b_f[1..|b_{f+1}|] \approx b_2b_3\ldots b_fb_{f+1}$, each position $x \in [1..\Delta]$ satisfies exactly one of the following possibilities:

1. $s[x + p \cdot \Delta]$ is the same, for all integers $p \in [0,f)$,
2. $s[x + p \cdot \Delta] < s[x + (p + 1) \cdot \Delta]$ for all integers $p \in [0,f-1)$,
3. $s[x + p \cdot \Delta] > s[x + (p + 1) \cdot \Delta]$ for all integers $p \in [0,f-1)$.

Note that $i_0 = i - \Delta$ satisfies (2) or (3), because $s[i]$ is different than $s[1], s[2], \ldots, s[i-1]$, so in particular $s[i-\Delta] \neq s[i]$.

By reversing the order of the alphabet, it is enough to establish the lemma assuming that $i_0$ satisfies (2). To this end, we choose some positions $i_1, i_2, \ldots, i_\ell$ in $b_1$ as follows. Let $C$ be the set of characters that appear in $b_1$. The position $i_1 \in [1,\Delta]$ is chosen so that $s[i_1]$ is the strict successor of $s[i_0]$ in $C$, then $i_2 \in [1,\Delta]$ is chosen so that $s[i_2]$ is the strict successor of $s[i_1]$ in $C$, and so on. If there are multiple choices for the next $i_j \in [1,\Delta]$ then we take the smallest. We stop when one of the following two possibilities holds:

(a) $s[i_\ell+1]$ is not defined, i.e. $s[i_\ell]$ is the largest character in $b_1$.

(b) $i_{\ell+1}$ satisfies (1) or (3).

Notice that, by definition, the positions $i_0, i_1, \ldots, i_\ell$ all satisfy (2). Further, $s[i_0]$ is a leftmost character because $s[i]$ is a leftmost character, so $s[i'] \neq s[i]$ for every $i' \in [1,i)$, and $b_1 \approx b_2$ so $s[i'-\Delta] \neq s[i-\Delta]$ for every $i' \in [\Delta+1,i)$. Next, we note that $s[i_1], s[i_2], \ldots, s[i_\ell]$ are all leftmost characters because we are always choosing the smallest $i_j$ such that $s[i_j]$ is equal to a specific character, for $j = 1, 2, \ldots, \ell$.

We summarize the situation so far. For every integer $p \in [0,f)$, the fragment $s[i_\ell + p \cdot \Delta]$ belongs to block $b_{p+1}$, and we want to show that it is a leftmost character. We know that $s[i_\ell]$ is a leftmost character, thus by $b_1 \approx b_{p+1}$ we obtain that $s[i_\ell + p \cdot \Delta]$ does not occur earlier in $b_{p+1}$. We need to establish that it also does not occur earlier in $b_1, b_2, \ldots, b_p$. We separately consider the two possible cases (a) and (b).
(a) $s[i_{\ell+1}]$ is not defined, i.e. $s[i_{\ell}]$ is the largest character in $b_1$. We know that $i_{\ell}$ satisfies (2), so $s[i_{\ell}] < s[i_{\ell} + |\Delta|] < \ldots < s[i_{\ell} + (p - 1) \cdot |\Delta|] < s[i_{\ell} + p \cdot |\Delta|]$. For all integers $q \in [0, p)$, by $b_1 \approx b_{q+1}$ we obtain that $s[i_{\ell} + q \cdot |\Delta|]$ is the largest character in $b_{q+1}$. So in fact $s[i_{\ell} + p \cdot |\Delta|]$ is larger than all characters in the whole block $b_{q+1}$, for every integer $q \in [0, p)$, making $i_{\ell} + p \cdot |\Delta|$ a leftmost occurrence.

(b) $i_{\ell+1}$ is defined and satisfies (1) or (3), so $s[i_{\ell+1}] \geq s[i_{\ell+1} + |\Delta|] \geq \ldots \geq s[i_{\ell+1} + (p-1) \cdot |\Delta|] \geq s[i_{\ell+1} + p \cdot |\Delta|]$. See Figure 8. We know that $i_{\ell}$ satisfies (2), so $s[i_{\ell}] < s[i_{\ell} + |\Delta|] < \ldots < s[i_{\ell} + (p - 1) \cdot |\Delta|] < s[i_{\ell} + p \cdot |\Delta|]$. Recall that $s[i_{\ell+1}]$ is a strict successor of $s[i_{\ell}]$ in $b_1$. Thus, for every $i^\prime \in [1, \Delta]$ we have that $s[i^\prime]$ does not belong to the interval $(s[i_{\ell}], s[i_{\ell+1}])$. Hence, for every integer $q \in [0, p)$, this implies $s[i^\prime + q \cdot |\Delta|]$ does not belong to the interval $(s[i_{\ell} + q \cdot |\Delta|], s[i_{\ell+1} + q \cdot |\Delta|])$. As observed earlier, $s[i_{\ell} + q \cdot |\Delta|] < s[i_{\ell} + p \cdot |\Delta|]$ and $s[i_{\ell+1} + q \cdot |\Delta|] \geq s[i_{\ell+1} + p \cdot |\Delta|]$. We conclude that, for every $i^\prime \in [1, \Delta]$, we have that $s[i^\prime + q \cdot |\Delta|]$ does not belong to the interval $[s[i_{\ell} + p \cdot |\Delta|], s[i_{\ell+1} + p \cdot |\Delta|])$ (the interval is non-empty, as both positions belong to the same block $b_q$, and by $b_{q+1} \approx b_1$ we have that $s[i_{\ell+1} + q \cdot |\Delta|]$ is a strict successor of $s[i_{\ell} + q \cdot |\Delta|]$ in $b_q$). In particular, $s[i^\prime + q \cdot |\Delta|] \neq s[i_{\ell} + p \cdot |\Delta|]$, so $s[i_{\ell} + p \cdot |\Delta|]$ does not occur in $b_{q+1}$, making it a leftmost character.

Hence, for every integer $p \in [0, f)$, position $i_{\ell} + p \cdot |\Delta|$ is a leftmost occurrence. ▶

![Figure 8](image-url) The red points correspond to $s[i_{\ell+1}], \ldots, s[i_{\ell+1} + p \cdot |\Delta|]$. The blue points correspond to $s[i_{\ell}], \ldots, s[i_{\ell} + p \cdot |\Delta|]$. The black arrow illustrates the character’s axis.

By combining the above lemmas we obtain the following conclusion.

Lemma 12. If $|O_k| \geq 3$ then for any $\{uw, u'v', u''v''\} \in O_k$ where $|u| < |u'| < |u''|$ such that $\{uw, u'v'\}$ defines $b_1, \ldots, b_f, b_{f+1}$ and $\{u'v', u''v''\}$ defines $b'_1, \ldots, b'_f, b'_{f+1}$ there is a leftmost occurrence in every block $b_1, b_2, \ldots, b_f$ or there is a leftmost occurrence in every block $b'_1, b'_2, \ldots, b'_f$.

Proof. Recall that by Proposition 7, $|u'| - |u|$ is an op-border of $u = s[1..|u|]$ while $|u''| - |u'|$ is an op-border of $u' = s[1..|u'|]$. By Lemma 10 there is a leftmost occurrence in block $b_2$ or in block $b'_2$. Then, by Lemma 11 applied either to the blocks $b_1, b_2, \ldots, b_f, b_{f+1}$ defined by the op-border $|u'| - |u|$ or the blocks $b'_1, b'_2, \ldots, b'_f, b'_{f+1}$ defined by the op-border $|u''| - |u'|$, there is a leftmost occurrence in every block $b_1, b_2, \ldots, b_f$ or in every block $b'_1, b'_2, \ldots, b'_f$. ▶

We are now ready to upper bound $\sum_k |O_k|$ by the number of leftmost characters. We will show that, if some $O_k$ is large then there are many leftmost characters in some range. To this end, we define groups of leftmost occurrences. Let $L_k$ be the set of the leftmost occurrences $i$ such that $i \in [2^k, 2^{k+1})$:.

Definition 13. $L_k = \{i \mid i \in [2^k, 2^{k+1}) \land \forall j \in [1, i), s[i] \neq s[j]\}$ for $0 \leq k \leq \log m$. 

Note that the groups are disjoint, i.e. $L_k \cap L_{k'} = \emptyset$ for any $k$ and $k'$. Thus $\sum_{k=0}^{\log m}|L_k| \leq \sigma$. With this definition in hand, we are ready to show the main technical lemma.

Lemma 14. The number of op-squares that are prefixes of $s$ is $O(\sigma)$.

Proof. To establish the lemma we want to connect $|O_k|$ with $|L_k|$, and then sum over all possible values of $k$. $k = 0, 1$ will be considered separately, and for larger $k$ we apply different arguments for $|O_k| \geq 11$ and $|O_k| \leq 10$.

We first consider $k \geq 2$ such that $|O_k| \geq 11$. In particular, when $|O_k| \geq 3$, so by Proposition 8, there exist $\{uv, u'v', u''v''\} \in O_k$ such that $0 < |u'| - |u|, |u''| - |u'| < 2^k/(|O_k| - 2)$. By Lemma 12, for any $\{uv, u'v', u''v''\} \in O_k$ where $|u| < |u'| < |u''|$ such that $\{uv, u'v'\}$ defines $b_1, \ldots, b_f, b_{f+1}$ in $u$ and $\{u'v', u''v''\}$ defines $b_1', \ldots, b_{f'}, b'_{f+1}$ in $u'$ either there is a leftmost occurrence in every block $b_1, b_2, \ldots, b_f$ or there is a leftmost occurrence in every block $b'_1, b'_2, \ldots, b'_{f'}$. In either case, we have found $\{uv, u'v'\} \in O_k$ with $0 < \Delta < 2^k/(|O_k| - 2)$, where $\Delta = |u'| - |u|$, such that $\{uv, u'v'\}$ defines $b_1, \ldots, b_f, b_{f+1}$ with $f = \left\lfloor \frac{|u|}{\Delta} \right\rfloor$ and there is a leftmost occurrence in every block $b_1, b_2, \ldots, b_f$. We want to establish a lower bound on the number of leftmost occurrences in $L_{k-2}$. To this end, it is enough to show a lower bound on the number of blocks $b_i$ that are fully contained in the range $[2^{k-2}, 2^{k-1})$. Recall that $|u| \in [2^{k-1}, 2^k)$, and $u = b_1b_2 \ldots b_{f+1}$. Thus, the fragment $u[2^{k-2}, 2^{k-1} - 1]$ consists of a suffix (possibly empty) of some $b_j$, then $b_{j+1}, b_{j+2}, \ldots, b_{j+\ell}$, and then a prefix of $b_{j+\ell+1}$ (where $b_{j+\ell+1}$ might be the incomplete block $b_{j+\ell}$ that should not be counted in the lower bound). Thus, the number of blocks $b_i$ that are fully contained in the range $[2^{k-2}, 2^{k-1})$ is at least $2^{k-2}/\Delta - 1$. See Figure 9. Combining this with the upper bound on $\Delta$, we obtain the following inequality:

$$|L_{k-2}| \geq \frac{2^{k-2}}{\Delta} - 1 \geq \frac{2^{k-2}}{\Delta} - 2 > \frac{|O_k| - 2}{4} - 2 = \frac{|O_k| - 10}{4}. $$

Using the assumption $|O_k| \geq 11$, we conclude that $|L_{k-2}| > |O_k|/44$. Hence:

$$\sum_{k=2}^{\log m} |O_k| < \sum_{k=2}^{10} 44 \cdot |L_{k-2}| \leq 44 \sum_k |L_k| \leq 44 \cdot \sigma.$$
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\[
\sum_{k \geq 2} |O_k| \leq \sum_{k \geq 2} |O_k| + 10 + \sum_{k \geq 2} |L_{k-1}| + \sum_{k \geq 2} |L_k| \leq 20\sigma.
\]

To upper bound \(\sum_k |O_k|\), we split the sum into three parts. For \(k = 0, 1\), we have \(|O_k| \leq 1\) and \(|O_1| \leq 2\). Then, for \(k \geq 2\) we separately consider all \(k\) with \(|O_k| \geq 11\) and \(|O_k| \leq 10\) and plug in the above upper bounds. Overall, we obtain:

\[
\sum_k |O_k| \leq 1 + 2 + 44 \cdot \sigma + 20 \cdot \sigma = \mathcal{O}(\sigma).
\]

Thus, the number of op-squares that are prefixes of \(s\) is \(\mathcal{O}(\sigma)\).

We conclude the section with the main theorem.

**Theorem 15.** The number of op-squares in a string \(w\) of length \(n\) over an alphabet of size \(\sigma\) is \(\mathcal{O}(\sigma n)\).

**Proof.** We consider each suffix of \(w\) separately. For each suffix \(w[i..n]\), we apply Lemma 14 to conclude that the number of op-squares that are prefixes of \(w[i..n]\) is upper bounded by \(\mathcal{O}(\sigma)\). Thus, summing over all \(i\) we obtain that the number of op-squares in \(w\) is \(\mathcal{O}(\sigma n)\).

5 Algorithm

In this section, we describe the algorithm that reports all occurrences of op-squares in a string \(w[1..n]\) over an alphabet of size \(\sigma\) in \(\mathcal{O}(\sigma n)\) time.

The high-level idea of the algorithm is to generate \(\mathcal{O}(\sigma n)\) candidates for op-squares and then test each of them in constant time, see the pseudocode in Algorithm 1. To this end, we first describe a mechanism for checking if \(w[i..i+\ell-1] \approx w[i+\ell..i+2\ell-1]\) in constant time. This can be implemented with an LCA query on the order-preserving suffix tree of \(w\), as explained in [3]. However, we need to explain how to construct this structure in \(\mathcal{O}(\sigma n)\) time.

**Order-preserving suffix tree.** Following [3], for a string \(w[1..n]\) we define code\((w)\) as \((\phi(w, 1), \phi(w, 2), \ldots, \phi(w, n))\), where \(\phi(w, i) = (\text{prev}_<(w, i), \text{prev}_>(w, i))\) and \(\text{prev}_<(w, i) = \{|k < i : w[k] < w[i]\}|, \text{prev}_>(w, i) = \{|k < i : w[k] = w[i]\}|\). We observe that code\((w)\) = code\((w')\) if and only if \(w \approx w'\). Then, the order-preserving suffix tree of \(w[1..n]\) is the compacted trie of all strings of the form code\((w[i..n])\)$, for \(i = 1, 2, \ldots, n\). It is easy to see that \(w[i..i+\ell-1] \approx w[i+\ell..i+2\ell-1]\) if and only if the lowest common ancestor of the leaves corresponding to code\((w[i..n])\)$ and code\((w[i+\ell..n])\)$ is at string depth at least \(\ell\). Therefore, assuming that we have already built the order-preserving suffix tree of \(w[1..n]\), such a test can be implemented in constant time after \(\mathcal{O}(n)\) preprocessing for LCA queries [15]. It remains to explain how to construct the order-preserving suffix tree. We stress that while [3] does provide an efficient \(\mathcal{O}(n \log n / \log \log n)\) time construction algorithm (in fact, the full version [4] further improves the time complexity to \(\mathcal{O}(n \sqrt{\log n})\)), such complexity is incompatible with our goal. Due to the lack of space, the proof is moved to the appendix.

**Lemma 16.** Given a string \(w[1..n]\) over an alphabet of size \(\sigma\), we can construct its order-preserving suffix tree in \(\mathcal{O}(\sigma n)\) time and space.

The main part of the algorithm is efficiently generating \(\mathcal{O}(\sigma n)\) candidates for op-squares. Then, each of them is tested in constant time as explained above, assuming the preprocessing from Lemma 16.
\begin{algorithm}
\caption{Report all occurrences of op-squares in a string \(w[1..n]\) over an alphabet of size \(\sigma\) in \(O(\sigma n)\) time.}
\begin{algorithmic}[1]
\State Preprocess \(w[1..n]\) for retrieving the characters of any \(\text{code}(w[i..n])\)
\State Construct the order-preserving suffix tree \(T\)
\State Preprocess \(T\) for LCA queries
\State \(i \leftarrow n\)
\While {\(i > 0\)}
\State \(s \leftarrow w[i..n]\)
\ForEach {\(x_j\) that is the smallest or the largest of its group \(L_k\) do}
\State \(s' \leftarrow s[1..2^{k-1}]\)
\State \(\text{foreach fragment } s[y..y + 2^{k-1} - 1] \approx s'\) such that \(x_j \in [y, y + 2^{k-1} - 1]\) do
\State \hspace{1em} Store \(s[1..2(y - 1)]\) as a candidate in \(R[x_j][k][i]\)
\EndFor
\EndFor
\ForEach {candidate \(s[1..2(y - 1)]\) in \(R[x_j][k][i]\) do}
\State \(v_1\) $\leftarrow$ the leaf corresponding to \(\text{code}(w[i..n])\) in \(T\)
\State \(v_2\) $\leftarrow$ the leaf corresponding to \(\text{code}(w[i + (y - 1)..n])\) in \(T\)
\If {the string depth of \(LCA_T(v_1, v_2)\) is at least \(y - 1\)}
\State Report \(s[1..2(y - 1) - 1]\) as an op-square
\EndIf
\EndFor
\State \(i \leftarrow i - 1\)
\EndWhile
\end{algorithmic}
\end{algorithm}

Leftmost occurrences. As in the proof of the \(O(\sigma n)\) upper bound on the number of op-squares, we will consider the suffixes of the input string \(w[1..n]\) one-by-one. For \(i = n, n - 1, \ldots, 1\) in this order, let \(s = w[i..n]\) be the currently considered suffix, and \(x_1 < x_2 < \ldots < x_t\) be the leftmost occurrences in \(s\). By spending \(O(\sigma)\) time per each suffix, we can assume that the positions \(x_1, x_2, \ldots, x_t\) are known, as after moving from \(w[i..n]\) to \(w[i - 1..n]\) we only have to insert the new leftmost occurrence \(i - 1\) and possibly remove the previous leftmost occurrence \(i'\) such that \(w[i - 1] = w[i']\) (unless \(w[i - 1]\) has not been seen before), which can be done in \(O(t) = O(\sigma)\) time. By Proposition 5, every prefix of \(s\) that is an op-square can be obtained by choosing two leftmost characters at positions \(x_q\) and \(x_j\), where \(q < j\), and setting the length of the possible square to be \(2(x_j - x_q)\). This gives us \(O(\sigma^2)\) candidates for prefixes that could be op-squares. However, our goal is to generate only \(O(\sigma)\) such candidates. To achieve this goal, we first provide some combinatorial properties in Lemma 17, Lemma 18, and Proposition 19.

Recall that all leftmost occurrences are partitioned into groups \(L_0, L_1, \ldots\). Next, we show that it is enough to consider \(x_j\) that is the smallest or the largest element in its group.

\begin{lemma}
Consider an op-square \(s[1..2\ell]\). Then there exists \(q < j\) such that the leftmost occurrences \(x_q\) and \(x_j\) satisfy \(x_j - x_q = \ell\), \(x_j \in [\ell + 1, 2\ell]\) and \(x_j\) is either the smallest or the largest element of its group.
\end{lemma}

The proof is described in the appendix. Figure 10 illustrates the scenario of the lemma.

To generate the candidates, we iterate over all \(j\) such that \(x_j\) is the smallest or largest element of its group \(L_k\). Consider \(q < j\) such that \(x_j - x_q = \ell\) and \(x_j \in [\ell + 1, 2\ell]\) for an op-square \(s[1..2\ell]\). Then, because \(x_j \in [2^k, 2^{k+1}]\), \(\ell \geq 2^{k-1}\). To avoid clutter, let \(s' = s[1..2^{k-1}]\).
Because \( s[1..2\ell] \) is assumed to be an op-square, we have \( s[\ell, \ell + 2^{k-1} - 1] \approx s' \). This suggests the following natural strategy to generate the candidates: we iterate over all fragments \( s[y..y + 2^{k-1} - 1] \) such that \( x_j \in [y, y + 2^{k-1} - 1] \) and \( s' \approx s[y..y + 2^{k-1} - 1] \), and output \( s[1..(y-1)] \) as a possible op-square (as explained earlier, each such candidate is then tested in constant time). See Figure 11. We first bound the number of such fragments by \( O(1 + |L_{k-2}|) \), and then explain how to generate them in the same time complexity.

### Lemma 18.
The number of fragments \( s[y..y + 2^{k-1} - 1] \) such that \( x_j \in [y, y + 2^{k-1} - 1] \) and \( s' \approx s[y..y + 2^{k-1} - 1] \) is upper bounded by \( O(1 + |L_{k-2}|) \).

The proof of the lemma relies on showing that \( \ell' \) is an op-border of \( s' \) and thus we can define blocks of length \( \Delta = 2^{k-1} - \ell' \) in \( s' \) and then apply Lemma 10 and Lemma 11 to achieve the desired bound. The full proof is described in the appendix. Hence, for every \( k \) such that \( L_k \) is non-empty, we generate \( O(1 + |L_{k-2}|) \) candidates. The overall number of candidates generated by following the above strategy is \( \sum_{k: L_k \neq \emptyset} O(1 + |L_{k-2}|) = O(\sigma + \sum_k |L_{k-2}|) = O(\sigma) \) as promised. It remains to show how to access all fragments \( s[y..y + 2^{k-1} - 1] \) such that \( x_j \in [y, y + 2^{k-1} - 1] \) and \( s' \approx s[y..y + 2^{k-1} - 1] \) in time proportional to their number.

### Accessing candidates.
We will solve a more general problem, and show how to ensure that, when considering \( s = w[i..n] \), for every leftmost occurrence \( x_j \) in \( s \) we have access to a list of all fragments \( s[y..y + 2^{k-1} - 1] \) such that \( x_j \in [y, y + 2^{k-1} - 1] \) and \( s[1..2^{k-1} - 1] \approx s[y..y + 2^{k-1} - 1] \), where \( 2^k \leq x_j < 2^{k+1} \). We call this list the result for \( i \) and \( x_j \).

Recall that \( s = w[i..n] \), and we consider \( i = n, n - 1, \ldots, 1 \) in this order. When we consider \( s = w[i..n] \), position \( i \) becomes a leftmost occurrence and remains to be so until we reach \( s = w[\text{prev}_i..n] \) such that \( w[i] = w[\text{prev}_i] \) (possibly, it is a leftmost occurrence till the very end of the scan). We can calculate \( \text{prev} \) for every \( i \) in \( O(\sigma n) \) time by maintaining a list of leftmost occurrences as described earlier. We say that a position \( i \) is \( k \)-active at position \( i' \) when \( i' \in [\text{prev}_i, i] \) and \( 2^k \leq i - i' + 1 < 2^{k+1} \). We observe that, as we consider longer and longer suffixes of \( w \), position \( i \) is first 0-active, then 1-active, and so on until it becomes...
We observe that every position $i'$ such that $i$ is $k$-active at $i'$ form a contiguous range $[\text{begin}_{i,k}, \text{end}_{i,k}]$ (the length of each such range is $2^k$, except possibly for $k = k_i$ when it is shorter). The total length of these ranges is small as shown below.

**Proposition 19.** $\sum_{i,k:k \leq k_i} 2^k = O(\sigma n)$

Proof. For $k \geq 1$ we can upper bound $2^k$ by $2 \cdot [\text{begin}_{i,k-1}, \text{end}_{i,k-1}]$. Then the sum becomes:

$$\sum_{i,k:k \leq k_i} 2^k = n + 2 \cdot \sum_{i,k:1 \leq k \leq k_i} [\text{begin}_{i,k-1}, \text{end}_{i,k-1}] \leq n + 2 \cdot \sum_{i,k:1 \leq k \leq k_i} [\text{begin}_{i,k}, \text{end}_{i,k}].$$

We observe that every position $i' \in [\text{begin}_{i,k}, \text{end}_{i,k}]$ in the suffix $w[i..n]$ corresponds to the relative position $i - i' + 1$ being a leftmost occurrence in the suffix $w[i'..n]$. Because there are at most $\sigma$ leftmost characters in any suffix $w[i'..n]$, this allows us to upper bound the sum by $O(\sigma n)$.

**Storing candidates.** This allows us to physically store the results as follows. For every leftmost occurrence $x$, we have an array indexed by $k \leq k_i$, denoted $R[x]$. Each entry of this array is an array indexed by $i' \in [\text{begin}_{i,k}, \text{end}_{i,k}]$, denoted $R[x][k][i']$. Finally, each entry of that array, denoted $R[x][k][i']$, is a pointer to a list of $y$s such that $x \in [y, y + 2^{k-1} - 1]$ and $w[i'..x'] + 2^{k-1} - 1 \approx w[y..y + 2^k - 1] - 1$ (note that it is a pointer to a list and not its physical copy). The arrays allow us to access the result for every $x, y$, $k \leq k_x$, and $i'$ in constant time, by retrieving the pointer $R[x][k][i']$ (where we first verify that $i' \in [\text{begin}_{i,k}, \text{end}_{i,k}]$). The total length of all arrays $R[x]$ is only $O(\sigma n)$ by Proposition 19. Further, the total length of all lists of occurrences that we need to prepare (assuming that we store every $R[x][k][i]$ as a pointer to such a list and not their physical copies) is also $O(\sigma n)$ by the following argument. Consider $i$ and $k \leq k_i$. Then, we need a list of positions $y$ such that $i \in [y, y + 2^k - 1]$ and $w[y..y + 2^{k-1} - 1]$ is order-isomorphic to a specific string $s'$. Thus, we can partition all positions $y$ such that $i \in [y, y + 2^k - 1]$ into groups corresponding to order-isomorphic fragments $w[y..y + 2^{k-1} - 1]$, and then store a pointer to the appropriate list (possibly null, if there is no $y$). The total number of positions $y$, over all $i$ and $k \leq k_i$, is $O(\sigma n)$ by Proposition 19, which bounds the total length of all the lists.

**Generating candidates.** It remains to describe how to efficiently calculate the results. This requires partitioning all fragments $w[y..y + 2^{k-1} - 1]$ such that $i \in [y, y + 2^k - 1]$ and $k \leq k_i$ into order-isomorphic groups, and finding for every $i, k \leq k_i, i' \in [\text{begin}_{i,k}, \text{end}_{i,k}]$ a pointer to the list of fragments $w[y..y + 2^{k-1} - 1]$ with $i \in [y, y + 2^k - 1]$ that are order-isomorphic to $w[i'..x'] + 2^{k-1} - 1$. Both steps can be implemented with the order-preserving suffix tree that is preprocessed in $O(n)$ time and space for computing a (deterministic) fingerprint of any code($w[x..x + 2^k - 1]$) in constant time. Here, a fingerprint is meant as an integer consisting of $O(\log n)$ bits, denoted fingerprint$(x)$, such that fingerprint$(x) = \text{fingerprint}(x')$ iff code($w[x..x + 2^k - 1]$) = code($w[x'..x' + 2^k - 1]$) (or equivalently $w[x..x + 2^k - 1] \approx w[x'..x' + 2^k - 1]$). We first describe such a mechanism and then provide a more detailed description of how to apply it. The proof of the following lemma directly follows from prior work [11,12] and is described in the appendix.

**Lemma 20.** A compacted trie on $n$ leaves can be preprocessed in $O(n)$ time, so that for any leaf $u$ and integer $k$ we can query in constant time for a $O(\log n)$-bit fingerprint of the ancestor of $u$ at string depth $2^k$. 

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We apply Lemma 20 on the order-preserving suffix tree. This allows us to calculate any fingerprint(x) with the required properties in constant time. Now consider any i and k ≤ k_i. We first compute fingerprint_k_{i−1}(y) for every y such that i ∈ [y, y + 2^k_i − 1]. This takes O(2^k) time. Next, we compute fingerprint_k_{i−1}(i') for every i' ∈ [begin_{i,k}, end_{i,k}], also in O(2^k) time because |[begin_{i,k}, end_{i,k}]| ≤ 2^{k−1}. We sort all fingerprints and partition them into groups corresponding to order-isomorphic fragments. We need to implement this step in O(2^{k}) time as well. To this end, we observe that we need to sort O(2^k) integers consisting of O(log n) bits, which can be done with radix sort in O(2^k + n) time. To avoid paying O(n) for each i and k ≤ k_i, we observe that this is an offline problem, and all sets corresponding to different i and k ≤ k_i can be sorted together. In more detail, we sort tuples of the form (i, k_i, fingerprint_y(y), y) and (i, k_i, fingerprint_y(i'), i'). The total number of all tuples is O(σn) by Lemma 19 and, as each of them can be treated as an integer consisting of O(log n) bits, they can be sorted in O(σn + n) = O(σn) time. Then, we extract the results for each i and k ≤ k_i from the output. For each i and k ≤ k_i, we consider every group of equal fingerprints. From each group, we first create a list containing all positions y corresponding to fingerprint_k_{i−1}(y) belonging to the group. Then, for every fingerprint_k_{i−1}(i') belonging to the group we store a pointer to this list. Overall, this takes O(σn) time and allows us to compute all the results in the same time complexity.

6 Open Problems

An interesting follow-up to our results is first bounding the number of order-preserving squares that are not order-isomorphic, and then designing an algorithm that reports all such squares. In addition, investigating the bounds for parameterized squares is of interest. Moreover, we are not aware of an algorithm reporting parameterized squares in a string, hence, designing such an algorithm is desired.

References

Lemma 16. Given a string \(w[1..n]\) over an alphabet of size \(\sigma\), we can construct its order-preserving suffix tree in \(O(\sigma n)\) time and space.

Proof. As explained in [3], the order-preserving suffix tree of \(w[1..n]\) can be constructed using the general framework of Cole and Hariharan [2] for constructing a suffix tree for a quasi-suffix collection of strings \(w_1, w_2, \ldots, w_n\). The running time of their algorithm is \(O(n)\) with almost inverse exponential failure probability, assuming that one can access the \(j\)-th character of any \(w_i\) in constant time. The mechanism for accessing the \(j\)-th character of \(w_i\) is called the character oracle. In this particular application, the string \(w_i = \phi(w[i..n])\). We will first describe how to implement a constant-time character oracle for such strings, and then explain why randomization is not needed in our setting.
We need to implement a new character oracle that returns $\phi(w[i..n], j)$, for any $i, j$, in constant time after $O(\sigma n)$ time and space preprocessing. This requires being able to calculate $\text{prev}_<(w[i..n], j)$ and $\text{prev}_>(w[i..n], j)$ in constant time. To this end, we define a two-dimensional array $\text{cnt}[i, x] = |\{k : k < i, w[k] < x\}|$, for $i = 0, 1, \ldots, n$ and $x = 0, 1, \ldots, \sigma$. All entries in this array can be computed in $O(\sigma n)$ total time and space. Then, we can calculate any $\text{prev}_<(w[i..n], j)$ and $\text{prev}_>(w[i..n], j)$ as follows:

$$
\text{prev}_<(w[i..n], j) = \text{cnt}[i + j - 2, w[i + j - 1]] - \text{cnt}[i - 1, w[i + j - 1]]
$$

$$
\text{prev}_>(w[i..n], j) = (\text{cnt}[i + j - 2, w[i + j - 1]] - \text{cnt}[i + j - 2, w[i + j - 1] - 1]) - (\text{cnt}[i - 1, w[i + j - 1]] - \text{cnt}[i - 1, w[i + j - 1] - 1]).
$$

To remove randomization, we observe that its only source in the algorithm of Cole and Hariharan is the need to maintain, for each explicit node of the current tree, a dictionary indexed by the next character on an outgoing edge. If we could show that there are at most $O(\sigma)$ such edges, then the dictionary could be implemented as a simple list, increasing the construction time to $O(\sigma n)$, which is within our claimed bound.

Consider a non-leaf node $v$ of the current tree. It corresponds to a proper prefix of some $\text{code}(w[i..n])^8$, which by the definition of $\text{code}(\cdot)$ is equal to $\text{code}(w[i..j])$, for some $j$. Let $c_1 < c_2 < \ldots < c_k$ be the distinct characters of $w[i..j]$, and denote by $\text{occ}_x$ the number of occurrences of $c_x$ in $w[i..j]$. Now consider an edge outgoing from $v$, and let $\text{code}(w[i'..j'+1])$ correspond to the first node (implicit or explicit) after $v$ we know that $\text{code}(w[i'..j']) = \text{code}(w[i..j])$, so the distinct characters of $w[i'..j']$ are $c'_1 < c'_2 < \ldots < c'_k$ with $\text{occ}_{c_x}$ being the number of occurrences of $c'_x$ in $w[i'..j']$. Then, we analyze the possible values of $\text{prev}_<(w[i'..j'+1], j' - i' + 2), \text{prev}_>(w[i'..j'+1], j' - i' + 2)$, that is, the first character on the considered edge. The first number is always equal to $\sum_{y=1}^{x} c_y$, for some $x \in [1, k + 1]$. Then, the second number is either $0$ or $\text{occ}_x$. Thus, overall we have only $2k \leq 2\sigma$ possible first characters, which bounds the degree of any $v$ by $O(\sigma)$.

\begin{lemma}
Consider an op-square $s[1..2\ell]$. Then there exists $q < j$ such that the leftmost occurrences $x_q$ and $x_j$ satisfy $x_j - x_q = \ell$, $x_j \in [\ell + 1, 2\ell]$ and $x_j$ is either the smallest or the largest element of its group.
\end{lemma}

\begin{proof}
By Proposition 5, we know that there is a leftmost character in $s[\ell + 1..2\ell]$. Choose the largest $k$ such that $2^k \leq \ell$ (so $2^{k+1} > \ell$). Consider two ranges $[2^k, 2^{k+1})$ and $[2^{k+1}, 2^{k+2})$ corresponding to groups $L_k$ and $L_{k+1}$, respectively. Because $2^k \leq \ell$ and $2^{k+1} > \ell$, we have $2^k < \ell + 1$, $2^{k+1} \in [\ell + 1, 2\ell]$ and $2\ell < 2^{k+2}$. Consequently, the fragment $s[\ell + 1..2\ell]$ can be represented as the concatenation of a suffix of $s[2^k..2^{k+1})$ and a prefix of $s[2^{k+1}..2^{k+2})$. The leftmost occurrence that falls within $s[\ell + 1..2\ell]$ belongs to the suffix or the prefix. See Figure 10. If it falls within the suffix, the largest element of $L_k$ belongs to $[\ell + 1, 2\ell]$. If it falls within the prefix, the smallest element of $L_{k+1}$ belongs to $[\ell + 1, 2\ell]$. Let $x_j \in [\ell + 1, 2\ell]$ be the corresponding leftmost occurrence. To complete the proof we need to establish that there exists $q < j$ such that $x_j - x_q = \ell$. The character $s[x_j]$ is distinct from all $s[1], s[2], \ldots, s[x_j - 1]$, and by $s[1..\ell] \approx s[\ell + 1..2\ell]$ we obtain that $s[x_j - \ell]$ is distinct from all $s[1], s[2], \ldots, s[x_j - \ell - 1]$. Thus, the position $x_j - \ell$ is a leftmost occurrence, hence $x_j - \ell = x_q$ for some $q < j$ as required.
\end{proof}

\begin{lemma}
The number of fragments $s[y..y + 2^{k-1} - 1]$ such that $x_j \in [y, y + 2^{k-1} - 1]$ and $s' \approx s[y..y + 2^{k-1} - 1]$ is upper bounded by $O(1 + |L_{k-2}|)$.
\end{lemma}
Proof. Consider all such fragments \( s[y_1..y_1 + 2^{k-1} - 1], s[y_2..y_2 + 2^{k-1} - 1], \ldots, s[y_t..y_t + 2^{k-1} - 1] \). Because \( x_j \in [y_z, y_z + 2^{k-1} - 1] \) for every \( z = 1, 2, \ldots, t \), either \( t = 1 \) or by the pigeonhole principle there exists \( z \) such that \( y_{z+1} - y_z < 2^{k-1}/(t-1) \). If \( t = 1 \) then we are done. Otherwise, let \( t' = [s[y_{z+1}..y_z + 2^{k-1} - 1]] \). By assumption, \( s' \approx s[y_{z+1}..y_z + 2^{k-1} - 1] \) and \( s' \approx s[y_{z+1}..y_z + 2^{k-1} - 1] \), so by the transitivity of \( \approx \) also \( s[y_z..y_z + 2^{k-1} - 1] \approx s[y_{z+1}..y_{z+1} + 2^{k-1} - 1] \). We conclude that \( s'[1..t'] \approx s'[y_{z+1} - y_z + 1..2^{k-1}] \), or in other words \( t' \) is an op-border of \( s' \). Let \( b_1, b_2, \ldots, b_f \) be the blocks defined by \( t' \) in \( s' = s[1..|s'|] = w[i..i + |s'| - 1] \), where each block is of length \( \Delta = 2^{k-1} - t' \). See Figure 12. Recall that \( x_j \) is a leftmost occurrence in \( s = w[i..n] \), and by the definition of \( y_z \) and \( y_{z+1} \) we have \( x_j \in [y_{z+1}, y_z + 2^{k-1} - 1] \). Then, by \( s'[1..t'] \approx s'[y_{z+1} - y_z + 1..2^{k-1}] \) we obtain that \( x_j - y_z + 1 \in [y_{z+1} - y_z + 1, 2^{k-1}] \) is also a leftmost occurrence in \( s = w[i..n] \). Hence, we have a leftmost occurrence in block \( b_j \), for some \( j \geq 2 \). This allows us to apply Lemma 10 and then Lemma 11 to conclude that there is a leftmost occurrence in every block \( b_1, b_2, \ldots, b_f \).

We calculate a lower bound on how many of these leftmost occurrences fall within the range \([2^{k-2}, 2^{k-1}]\):

\[
\left\lfloor \frac{2^{k-2}}{\Delta} \right\rfloor - 1 > \frac{2^{k-2}}{2^{k-1} - t'} - 2
= \frac{2^{k-2}}{2^{k-1} - (y_z + 2^{k-1} - y_{z+1})} - 2
= \frac{2^{k-2} - 2}{y_{z+1} - y_z} - 2
> \frac{2^{k-2}}{2^{k-1}/(t-1)} - 2 = (t - 5)/2.
\]

For \( t < 6 \), we are done as the number of fragments is \( O(1) \). Otherwise, we obtain that \( |L_{k-2}| \geq (t - 5)/2 \geq t/12 \), thus \( t = O(1 + |L_{k-2}|) \) always holds as claimed.

\[\text{Figure 12} \ s, s', s[y_z..y_z + 2^{k-1} - 1], \text{and } s[y_{z+1}..y_{z+1} + 2^{k-1} - 1]\] are colored in black, orange, blue, and green, respectively. The red line illustrates the ranges.

\[\text{Lemma 20.} \ A \text{ compacted trie on } n \text{ leaves can be preprocessed in } O(n) \text{ time, so that for any leaf } u \text{ and integer } k \text{ we can query in constant time for a } O(\log n) \text{-bit fingerprint of the ancestor of } u \text{ at string depth } 2^k.\]

Proof. This follows by applying the method used to solve the substring fingerprint problem mentioned in [11, Lemma 14]. Following the description in the full version [12, Lemma 12], a compacted trie on \( n \) leaves can be preprocessed in \( O(n) \) time so that we can locate the (implicit or explicit) node corresponding to the ancestor at string depth \( 2^k \) of a given leaf in constant time. If the sought node is implicit (and does not explicitly exist in the compacted trie) we retrieve the edge that contains it. Next, if the node is explicit then we return its identifier. If the node is implicit then we return the identifier of the edge that contains it. Thus, the required range of identifiers is \([2n]\).
MUL-Tree Pruning for Consistency and Compatibility

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Abstract

A multi-labelled tree (or MUL-tree) is a rooted tree leaf-labelled by a set of labels, where each label may appear more than once in the tree. We consider the MUL-tree Set Pruning for Consistency problem (MULSETPC), which takes as input a set of MUL-trees and asks whether there exists a perfect pruning of each MUL-tree that results in a consistent set of single-labelled trees. MULSETPC was proven to be NP-complete by Gascon et al. when the MUL-trees are binary, each leaf label is used at most three times, and the number of MUL-trees is unbounded. To determine the computational complexity of the problem when the number of MUL-trees is constant was left as an open problem.

Here, we resolve this question by proving a much stronger result, namely that MULSETPC is NP-complete even when there are only two MUL-trees, every leaf label is used at most twice, and every MUL-tree is either binary or has constant height. Furthermore, we introduce an extension of MULSETPC that we call MULSETPC Comp, which replaces the notion of consistency with compatibility, and prove that MULSETPC Comp is NP-complete even when there are only two MUL-trees, every leaf label is used at most thrice, and every MUL-tree has constant height. Finally, we present a polynomial-time algorithm for instances of MULSETPC with a constant number of binary MUL-trees, in the special case where every leaf label occurs exactly once in at least one MUL-tree.

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1 Introduction

In evolutionary biology, leaf-labelled (phylogenetic) trees are commonly employed to describe the evolution of species using leaf labels to represent different species [11]. Comparisons of these structures are used particularly in phylogenetic inferences – similarities may indicate
evolutionary patterns, whereas differences may highlight genetic mutations. The measure of similarity between phylogenetic trees has been defined by multiple alternate metrics, such as the Robinson-Foulds distance [29], subtree pruning and regraft (SPR) distances [5, 34], and maximum agreement subtrees [2, 7, 12]. Other problems related to phylogenetic trees include constructing supertrees [1, 3, 4, 33] or consensus trees [6, 11, 21] which can determine relations or interactions between smaller phylogenetic trees.

Phylogenetic trees are classically described as single-labelled trees, where no label appears on the leaves of the tree more than once. Typically, construction or comparison algorithms of such phylogenetic trees make use of this property to reduce computational costs. Multi-labelled trees (or MUL-trees) are a generalisation of single-labelled trees in which multiple leaves may be labelled by the same label. MUL-trees can be useful to depict genome duplication, lineage sorting, or lateral gene transfer [23]. Other applications include the construction of phylogenetic networks by folding operations [17, 18, 19], biogeography [13, 24, 25], the study of host-parasite cospeciation [26], and gene evolution studies [23, 27, 30].

MUL-trees have been far less investigated than their single-labelled counterparts and many computational problems become NP-hard when extended to MUL-trees. For example, the majority rule consensus tree for a set of $k$ single-labelled trees with $n$ leaf labels each can be computed in $O(nk)$ time [21], but is NP-hard to compute for MUL-trees [8]. Other approaches convert MUL-trees into single-labelled trees which can be input to existing algorithms [20, 30]. A few polynomial-time algorithms do exist for MUL-trees – Cui et al. [8] presented a $O(n^2k + nk^2)$-time algorithm for building a majority rule consensus MUL-tree in which each leaf label occurs at most twice, based on a reduction to the Perfect Phylogeny Haplotyping problem [10]. Furthermore, the maximum agreement subtree (MAST) distance between two MUL-trees can be computed in quadratic time, though it also becomes NP-complete when generalised to more than two MUL-trees [13, 22].

This paper investigates MUL-trees by considering the MUL-tree Set Pruning for Consistency problem (MULSETPC), which takes as input a set of MUL-trees, and outputs whether or not there exists a pruning of the MUL-trees which gives a consistent set of single-labelled trees (see Section 2 for formal definitions). Gascon et al. showed that, in general, MULSETPC is NP-complete via a polynomial reduction from 3-SAT [15, 16]. However, their reduction from an instance of 3-SAT with $m$ variables and $z$ clauses gives an instance of MULSETPC containing $m + z + 1$ MUL-trees; moreover these MUL-trees may have labels occurring three times. Here we prove that MULSETPC is still NP-complete, even when restricted to instances involving only two MUL-trees, or in which every leaf label appears at most twice within a MUL-tree. This totally resolves the open question of Gascon et al. [15, 16] regarding the parameterised complexity of MULSETPC when the parameter is the number of input trees. Also, we identify tractable fragments of MULSETPC which can be solved in polynomial time, in short, instances in which each label appears exactly once in at least one MUL-tree.

We also present a generalisation of MULSETPC called MALSETPComp, which asks for a compatible set of trees instead of a consistent set of trees. Tree compatibility is a generalisation of tree consistency where we allow the supertree displaying the set to instead display a refinement of each tree rather than the tree itself—again, a more rigorous definition is given later. Tree compatibility is relevant when determining the existence of a supertree for a given set of phylogenetic trees [1, 31]. This is because experimental data often contains uncertainty, which can be expressed by non-binary nodes in the tree: this necessitates the use of compatibility. Compatibility is also relevant to other questions such as the incomplete directed perfect phylogeny problem [28]. Our interest in tree compatibility was partially motivated by recent improvements in compatibility testing [9]. Here, we prove that MALSETPComp is NP-complete, even when restricted to instances involving only two MUL-trees, or in which every leaf label appears at most thrice within a MUL-tree.
The rest of the paper is organized as follows. In Section 2, we introduce the preliminary notation and definitions. In Section 3 we present the improved NP-completeness proof for MULSETPC by reduction from the Boolean 3-SAT problem. In Section 4 we give a NP-completeness proof for MULSETPCmp by reduction from the Exact 3-cover with Multiplicity 3 problem. Section 5 contains our polynomial time results for tractable instances of MULSETPC. Finally, in Section 6 we present our conclusions and a few open problems.

2 Preliminaries

We shall use the following standard definitions on trees.

Definition 1 (Basic tree definitions). All trees we consider are rooted and unordered. If \( x, y \) are nodes in a tree \( T \), then \( y \) is an ancestor of \( x \) (and \( x \) a descendant of \( y \)) if \( y \) lies on the unique path from \( x \) to the root of \( T \). We denote this by \( x \leq y \). Additionally, if \( y \neq x \) then \( y \) is a proper ancestor of \( x \), which we denote by \( x < y \). If \( x < y \) and \( y \) is adjacent to \( x \) then \( y \) is the parent of \( x \) and \( x \) a child of \( y \). If \( x \) and \( x' \) are both children of \( y \) then \( x \) and \( x' \) are siblings. The lowest common ancestor of nodes \( x \) and \( y \), denoted \( \text{lca}_T(x, y) \), is the node \( z \) such that \( x \leq z \), \( y \leq z \), and no proper descendant of \( z \) also satisfies these properties. The empty tree, denoted by \( T_0 \), is the unique tree which contains no nodes.

We use the following definition of leaf-labelled trees, which takes the definitions of Gascon et. al. [15] and generalises them to the case where the tree may not be binary.

Definition 2 (Leaf-labelled trees). A leaf-labelled tree \( (T, \mathcal{X}) \) is a (rooted, unordered) tree \( T \) where no node has exactly one child and where each leaf has been assigned a label from a set of labels \( \mathcal{X} \). (We will sometimes refer to \( T \) as a leaf-labelled tree on \( \mathcal{X} \), and omit \( \mathcal{X} \) if it is clear from context.) A leaf-labelled tree is a single-labelled tree if every label in \( \mathcal{X} \) is used at most once. Alternatively, a multi-labelled tree or MUL-tree is a leaf-labelled tree where we allow each label in \( \mathcal{X} \) to label multiple leaves. We say a MUL-tree has multiplicity \( k \) if each leaf label appears at most \( k \) times. Let \( \mathcal{L}(T) \subseteq \mathcal{X} \) denote the set of leaf labels appearing in \( T \) and let \( D(T) \subseteq \mathcal{L}(T) \) denote the set of leaf labels appearing only once in \( T \).

Note that in a single-labelled tree we sometimes abuse notation and identify the leaf and the leaf label. We do the same in a MUL-tree only if the context is clear and there is no possibility of confusion.

If \( u \) is a node of \( T \) then \( T^u \) denotes the subtree rooted at \( u \) containing \( u \) and all its descendants, maintaining the same leaf-labelling as \( T \) on the remaining leaves. Let \( D^u := D(T^u) \).

Definition 3 (Pruning and Perfect Pruning). Given a leaf-labelled tree \( T \), let \( y \) denote a leaf node of \( T \) and \( x \) the parent of \( y \). We prune the leaf \( y \) in the following manner:

- Delete the leaf \( y \).
- If \( x \) still has at least two children, do nothing else.
- Alternatively, if \( x \) now has only one child and is not the root, suppress the vertex \( x \).
- Finally, if \( x \) has only one child \( z \) and is the root, delete \( x \) and make \( z \) the new root.

A perfect pruning of \( T \) is a single-labelled tree \( T' \) such that \( \mathcal{L}(T') = \mathcal{L}(T) \), created by (possibly repeated) prunings of \( T \). That is, for every label that appears more than once in \( T \), we prune away all but exactly one copy of the label to obtain a single-labelled tree. If \( T \) is a single-labelled tree, then its only perfect pruning is itself.
Given a leaf-labelled tree $T$ and a set of leaf labels $L' \subseteq L(T)$, let $T|_{L'}$ denote the leaf-labelled tree constructed by pruning from $T$ every leaf labelled by a label from $L(T) \setminus L'$. (That is, only leaves labelled by $L'$ remain.) Two leaf-labelled trees $T_1$ and $T_2$ are leaf-label isomorphic if there is an isomorphism between $T_1$ and $T_2$ which preserves the labelling of the leaves. We say that a leaf-labelled tree $T$ on $L$ displays a single-labelled tree $T'$ on $L' \subseteq L$ if there exists a perfect pruning $T^*$ of $T$ such that $T^*|_{L'}$ is leaf-label isomorphic to $T'$.

**Definition 4 (Refinement).** Given single-labelled trees $T$ and $T^*$, we say $T^*$ is a refinement of $T$ if $T$ can be obtained from $T^*$ by (possibly repeated) contractions of non-leaf edges, where we treat a contraction as merging the child node into the parent node. We write $T \preceq T^*$.

What follows is the definition of a **consistent set**, and the very similar definition of a **compatible set**.

**Definition 5 (Consistent Set).** Consider a set of single-labelled trees $T_1, \ldots, T_k$ with corresponding label sets $L_1, \ldots, L_k$. We say this set is **consistent** if there exists a single-labelled tree $T$ on label set $L = \bigcup_{i=1}^k L_i$ such that for every $i = 1, \ldots, k$, $T$ displays $T_i$.

Note in the above definition that if $L_1 = \cdots = L_k$ then $L = L_1$ and so the set $T_1, \ldots, T_k$ is consistent if and only if the trees are pairwise leaf-label isomorphic.

**Definition 6 (Compatible Set).** Consider a set of single-labelled trees $T_1, \ldots, T_k$ with corresponding label sets $L_1, \ldots, L_k$. We say this set is **compatible** if there exists a single-labelled tree $T$ on label set $L = \bigcup_{i=1}^k L_i$ such that for every $i = 1, \ldots, k$, $T|_{L_i}$ is a refinement of $T_i$.

![Figure 1](image)

Figure 1 Let $T_1$, $T_2$, and $T^*$ be the three trees with $\mathcal{X} = \{x, y, z, w\}$ shown above. If we prune away the non-underlined labels in $T_1$ and $T_2$ then $T^*|_{L(T_1)}$ is a refinement of the pruned $T_i$ for $i \in \{1, 2\}$, which shows that there exists a perfect pruning of $\{T_1, T_2\}$ giving a compatible set of trees. In contrast, there is no perfect pruning of $\{T_1, T_2\}$ giving a consistent set of trees because neither of the two single-labelled subtrees with leaf labels $y, z$, and $w$ displayed by $T_1$ is also displayed by $T_2$. Note, however, that if the label $w$ in $T_1$ is changed to $x$ then $\{T_1, T_2\}$ becomes consistent since this label can be pruned along with one leaf labelled by $z$ to obtain a perfect pruning of $T_1$ which is displayed by $T^*$.

Note that every consistent set of trees is also a compatible set, but the converse does not hold in general.

The MUL-set pruning for consistency problem can then be defined as follows:

<table>
<thead>
<tr>
<th>MUL-tree Set Pruning for Consistency (MULSETPC) Problem:</th>
</tr>
</thead>
<tbody>
<tr>
<td><strong>Input:</strong> $(\mathcal{M}, \mathcal{X})$ where $\mathcal{M}$ is a set of MUL-trees on $\mathcal{X}$.</td>
</tr>
<tr>
<td><strong>Output:</strong> $\exists? T$ a perfect pruning of each tree of $\mathcal{M}$ resulting in a consistent set of trees.</td>
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</table>
We introduce the following problem, which substitutes compatible for consistent sets:

**MUL-tree Set Pruning for Compatibility (MULSETPComp) Problem:**

**Input:** \((\mathcal{M}, \mathcal{X})\) where \(\mathcal{M}\) is a set of MUL-trees on \(\mathcal{X}\).

**Output:** \(\exists \gamma\) a perfect pruning of each tree of \(\mathcal{M}\) resulting in a compatible set of trees.

See Figure 1 for an example that illustrates the difference between MULSETPC and MULSETPComp.

### 3 NP-completeness for MULSETPC instances with two MUL-trees and multiplicity 2

In this section we consider the MULSETPC problem, and show that it is NP-complete even when considering a heavily restricted set of instances. Specifically, we consider instances with at most two MUL-trees, where the multiplicity is 2, and where the MUL-trees are binary. The core of our proof will be a reduction from 3-SAT [14].

**3-satisfiability (3-SAT) Problem:**

**Input:** A Boolean set of clauses \( C = (C_1 \land C_2 \land \ldots \land C_z) \) on a finite set of literals \( \{l_1, l_2, \ldots, l_m\} \) where each clause is in conjunctive normal and contains 3 literals.

**Output:** \( \exists \gamma\) a satisfying valuation \( V \) of \( C \).

Our first goal is to construct, given an instance of 3-SAT, two MUL-trees \( T_1 \) and \( T_2 \) which we will use in our corresponding instance of MULSETPC. Our set of leaf labels \( \mathcal{X} \) consists of the following:

\[
\{l_i, \bar{l}_i \mid i = 1, \ldots, m\}, \text{ the set of literals,}
\]

\[
\{F_i, V_i \mid i = 1, \ldots, m\}, \text{ a pair of “dummy” labels for each variable and}
\]

\[
\mathcal{P} := \{C_j^1, C_j^2, C_j^3 \mid j = 1, \ldots, z\}, \text{ a triple of “position” labels representing the three places of each clause of } C.
\]

Define \( h := \mathcal{P} \to \{l_i, \bar{l}_i \mid i = 1, \ldots, m\} \) as the function which maps a position label \( C_j^x \) to the literal found in that position. For example, if \( C_1 = (l_1 \lor l_4 \lor \bar{l}_6) \) then \( h(C_1^1) = l_1, h(C_1^2) = l_4 \) and \( h(C_1^3) = \bar{l}_6 \). We treat \( \mathcal{P} \) as being ordered first by the index of the clause and then by the position. Let \( H(l_i) := \{C_j^x \mid h(C_j^x) = l_i\} \), and define \( H(\bar{l}_i) \) similarly.

Our trees \( T_1, T_2 \) will be constructed from four types of subtrees \( T_{1,L,j}, T_{2,L,j}, T_{1,R,i}^*, T_{2,R,i} \), where \( i = 1, \ldots, m \) and \( j = 1, \ldots, z \). See Figure 2 for the subtrees \( T_{1,L,j}, T_{2,L,j}, T_{1,R,i}^* \) and \( T_{2,R,i} \); we now explain how to construct \( T_{1,R,i}^* \) and \( T_{2,R,i} \) from \( T_{1,R,i} \) and \( T_{2,R,i} \).

Let \( v_i \) denote the leaf labelled \( V_i \) in \( T_{1,R,i} \) and let \( u_i \) denote the parent of \( v_i \). Let \( u_i^+ \) and \( u_i^- \) denote the parents of leaves labelled \( l_i \) and \( \bar{l}_i \) respectively in \( T_{2,R,i} \). Let \( v_i^+ \) denote the child of \( u_i^+ \) labelled by \( V_i \), and \( v_i^- \) denote the child of \( u_i^- \) labelled by \( V_i \).
Initialise $T_{1,R,i}^*$ as a copy of $T_{1,R,i}$, and then:

- If $H(l_i) \cup H(\bar{l}_i) = \emptyset$, make no further changes.
- Otherwise, subdivide the edge $u_i v_i$ $|H(l_i) \cup H(\bar{l}_i)|$ times, and add to each new node an adjacent leaf. Label these leaves with $H(l_i) \cup H(\bar{l}_i)$, respecting the ordering such that the first label is closest to $u_i$.

Initialise $T_{2,R,i}^*$ as a copy of $T_{2,R,i}$, and then:

- If $H(l_i) \neq \emptyset$, then subdivide $u_i^+ v_i^- |H(l_i)|$ times and add an leaf adjacent to each new node. Label these leaves with $H(l_i)$, respecting the ordering so that the first label is closest to $u_i^+$.
- If $H(\bar{l}_i) \neq \emptyset$, repeat the previous step, substituting $H(\bar{l}_i)$ for $H(l_i)$ and $u_i^- v_i^-$ for $u_i^+ v_i^+$.

Construct $T_{1,L}$ by taking a complete binary tree on $z$ leaves (recall $z$ is the number of clauses), suppressing any nodes with exactly one child, and then identifying the root of each $T_{1,L,j}$ (ordered by $j$) with exactly one of the leaves (ordered left-to-right). Construct $T_{2,L}$ in the same fashion, substituting $T_{2,L,j}$ for $T_{1,L,j}$. Construct $T_{1,R}$ by taking a complete binary tree on $m$ leaves, suppressing any nodes with exactly one child, and then identifying the root of each $T_{1,R,i}^*$ (ordered by $i$) with exactly one of the leaves (ordered left-to-right). Again, construct $T_{2,R}$ in the same fashion, substituting $T_{2,R,i}^*$ for $T_{1,R,i}^*$. Finally, construct $T_1$ by taking a root node $r(T_1)$ and adding an edge to the roots of $T_{1,L}$ and $T_{1,R}$; construct $T_2$ in the obvious equivalent fashion.

Given an instance $C$ of 3-SAT, we create our instance of MULSETPC, $\langle \{T_1, T_2\}, X \rangle$. Note that $T_1$ and $T_2$ have multiplicity 2; each leaf label $C_j^x \in P$ appears once in $T_{1,L,j}$ and $T_{2,L,j}$ and once in $T_{1,R,i}$ and $T_{2,R,i}$ for the single value of $i$ such that $C_j^x \in H(l_i) \cup H(\bar{l}_i)$. By inspection, the labels of $X \sim P$ also appear at most twice. Hence the instance $\langle \{T_1, T_2\}, X \rangle$ contains two binary MUL-trees with multiplicity 2. It suffices to now show the reduction, in two parts.

**Lemma 7.** If $C$ is a satisfied instance of 3-SAT then the corresponding instance $\langle \{T_1, T_2\}, X \rangle$ of MULSETPC admits a perfect pruning giving a consistent set of trees.
Proof. Suppose that \( C \) is satisfiable. Then there exists a valuation of every variable which satisfies every clause of \( C \); fix one such valuation and label it \( \mathcal{V} \). For each clause \( C_j \), mark one of the position labels \( C_j^x \) for \( x \in \{1, 2, 3\} \) such that \( h(C_j^x) \) is valued true by \( \mathcal{V} \). Since \( \mathcal{V} \) satisfies every clause, we will always be able to choose a label to mark; if there are multiple legitimate choices choose arbitrarily. Refer to any label \( C_j^x \) we have not marked as unmarked.

By our construction of \( T_1 \) and \( T_2 \), if we show that after pruning each \( T_{1,L,j} \) and \( T_{1,R,i}^* \) is leaf-label isomorphic to \( T_{2,L,j} \) and \( T_{2,R,i}^* \) respectively, then \( T_1 \) is leaf-label isomorphic to \( T_2 \).

Consider first the labels of \( \mathcal{P} \). Prune from each \( T_{1,L,j} \) and \( T_{2,L,j} \) the one marked label \( C_j^x \), and leave the two unmarked labels \( C_j^y \). We must keep the other copy of the marked \( C_j^x \) and prune away the other copies of the unmarked \( C_j^y \) in whichever \( T_{1,R,i}^* \) and \( T_{2,R,i}^* \) they appear. Consider the sets \( H(l_i) \) and \( H(\bar{l}_i) \). If \( C_j^x \) is marked, then \( h(C_j^x) \) is true, and so at most one of \( H(l_i), H(\bar{l}_i) \) contains a marked label. Keeping this information in mind, we can now simply look at the subtrees themselves.

- After this pruning each \( T_{1,L,j} \) will be leaf-label isomorphic to the corresponding \( T_{2,L,j} \), by inspection.
- Consider \( T_{1,R,i}^* \) and \( T_{2,R,i}^* \). We have already pruned away unmarked labels of \( \mathcal{P} \). If \( l_i \) is true, prune the copy of \( l_i \) in \( T_{1,R,i}^* \) closest to the root of \( T_{1,R,i}^* \) and the copy of \( \bar{l}_i \) furthest from the root; if \( \bar{l}_i \) is true do the opposite. In \( T_{2,R,i}^* \) prune the copies of \( F_i, V_i \) closer to the literal \( l_i \) or \( \bar{l}_i \) which evaluates as false. Hence we have pruned away the extra copies of each leaf label. It is clear, mostly by inspection, that \( T_{1,R,i}^* \) is leaf-label isomorphic to \( T_{2,R,i}^* \); the most important point is that since at most one of \( H(l_i) \) and \( H(\bar{l}_i) \) contains a marked label, at least one of these sets has been pruned away entirely in \( T_{2,R,i}^* \) (specifically the set closer to the false literal).

Hence, after pruning, \( T_1 \) and \( T_2 \) are leaf-label isomorphic, and thus \( \{T_1, T_2\} \) is consistent. ▶

See Figure 3 for an illustration of the reduction in Lemma 7.

Lemma 8. If \( C \) is not a satisfied instance of 3-SAT then the corresponding instance \( \{T_1, T_2, X\} \) of \textsc{MULSETPC} do not admit perfect prunings giving a consistent set of trees.

We omit the proof of Lemma 8 for reasons of space. We now prove our main result.

Theorem 9. The \textsc{MULSETPC} problem is NP-complete, even restricted to instances containing at most two binary MUL-trees with multiplicity 2.

Proof. Note first that \textsc{MULSETPC} is in NP, since given a set of pruned leaves, a perfect pruning can be constructed in polynomial time, and the consistency of the set of trees determined in polynomial time using the BUILD algorithm [1, 15].

The result then follows directly from Lemma 7 and Lemma 8. ▶

It is also possible to swap our requirement that the MUL-trees are binary for an alternative requirement that the MUL-trees have height at most 5. Proving this result is very similar to the binary case, but we omit it here on grounds of space; the proof will appear in the journal version of this article. Thus we get the following result.

Theorem 10. The \textsc{MULSETPC} problem is NP-complete, even restricted to instances containing at most two MUL-trees with multiplicity 2 and height at most 5.
Figure 3 Illustrating the reduction from 3-SAT in Lemma 7. The two MUL-trees $T_1$ and $T_2$ are constructed from $C = (C_1 \land C_2 \land C_3)$ where $C_1 = (l_1 \lor l_2 \lor l_3)$, $C_2 = (l_2 \lor l_3 \lor l_4)$, and $C_3 = (l_1 \lor l_2 \lor l_4)$. Figure 3c shows the corresponding tree $T$ which displays the pruned $T_1$ and $T_2$ corresponding to a satisfiable assignment $l_1 = l_2 = l_3 = l_4 = \text{true}$ with marked labels $C_{c1}^b$, $C_{c2}^b$ and $C_{a3}^c$. 

(a) The MUL-tree $T_1$.

(b) The MUL-tree $T_2$.

(c) Tree $T$ displays the two MUL-trees $T_1$ and $T_2$ after they have been perfectly pruned.
4 NP-completeness for MULSETPCOMP

In this section, we extend our previous results regarding the problem MULSETPC to the similar problem MULSETPCOMP. We present the following two theorems.

▶ Theorem 11. MULSETPCOMP is NP-complete, even when restricted to instances containing at most two MUL-trees with multiplicity at most 3 and where the MUL-trees have height at most 4.

▶ Theorem 12. MULSETPCOMP is NP-complete, even when restricted to instances containing at most two MUL-trees with height at most 3 and where one MUL-tree is a single-labelled tree containing all leaf labels.

As was the case in Section 3, these two results have very similar proofs. We shall prove Theorem 11, but omit the proof of Theorem 12 for space reasons. Here, we reduce from X3C3, also known to be NP-complete [14].

Exact 3-cover with multiplicity 3 (X3C3) Problem:

Input: A set \( X = \{x_1, \ldots, x_{3q}\} \) and a collection \( C = \{S_1, \ldots, S_k\} \) of 3-element subsets of \( X \), such that any element of \( X \) appears in at most three sets in \( C \).

Output: \( \exists \? \) an exact cover for \( X \).

As before, our first goal is to construct, given an instance \((X, C)\) of X3C3, a set of two MUL-trees we shall use to construct our corresponding instance of MULSETPCOMP.

Recall that \(|X| = 3q\) and that \( k := |C| \). Let \( m := k - q \), the number of sets of \( C \) not chosen to be part of our exact 3 cover. Define \( A := \{a_{ij}^*|i = 1, \ldots, m, j = 1, \ldots, k\} \) and \( B = \{b_{ij}|i = 1, \ldots, m, j = 1, \ldots, k\} \). Let \( A^i = \{a_{ij}^*|j = 1, \ldots, k\} \) and \( A_j = \{a_{ij}^*|i = 1, \ldots, m\} \), and define \( B^i, B_j \) similarly. The labels in \( A \cup B \) are another set of “dummy” labels we use for technical reasons. Let \( Y = X \cup A \cup B \). We may assume that \( q \geq 3 \). We may also assume that \( k \) is even; if not, add to \( X \) three additional elements \( \{x_{3q+1}, x_{3q+2}, x_{3q+3}\} \) (which increases \( q \) by 1) and add to \( C \) a additional set \( S' = \{x_{3q+1}, x_{3q+2}, x_{3q+3}\} \) (which increases \( k = |C| \) by 1). It is clear this modified instance contains an exact 3-cover if and only if the original instance did.

We denote our two trees \( T_1 \) and \( T_2 \). Our corresponding instance of MULSETPCOMP for Theorem 11 will be \((T_1, T_2) \cup Y\).

\[ \text{Figure 4 The tree } T_1 \text{ for MULSETPCOMP.} \]
See Figures 4 and 6 for the construction of $T_1$, and Figures 5 and 7 for the construction of $T_2$. Note that a node labelled $\alpha^i$ in Figure 4 is the root of the appropriate subtree from Figure 6, not a leaf labelled by $\alpha^i$. An equivalent statement holds for $\bar{\alpha}^i, \beta^i$ and $\bar{\beta}^i$, where the subtree rooted at $\beta^i$ is found by taking the subtree of Figure 6 and replacing each $a^i_j$ with $b^i_j$. (An equivalent statement holds for $\bar{\beta}^i$ and Figure 7). Note the following other useful facts about our construction:

- Trees $T_1$ and $T_2$ have the same set of leaf labels. Hence a perfect pruning of $T_1$ and $T_2$ is a compatible set of trees if and only if there exists a tree $T^*$ on the same leaf label set which is a refinement of both perfect prunings.

- In $T_1$ any label of $X$ may appear at most three times, since any $x_i$ may appear in at most three sets of $C$. Every label of $A$ and $B$ appears once in $T^\mu_1$ and once in $T^\mu_2$. Hence $T_1$ has multiplicity 3. In $T_2$ any label of $X$ appears only once, and the labels of $A$ and $B$ appear at most thrice, once in $T^\nu_2$ and once or twice in $T^\nu_2$. Hence $T_2$ has multiplicity 3.

- The heights of both MUL-trees can be determined by inspection.

We will need the following two technical lemmas – as the proofs are straightforward we omit them.

▶ Lemma 13. Consider a set of elements $X = \{x_1, \ldots, x_k\}$ together with a subset $X' \subset X$ such that $|X'| = k - 1$ and a collection of sets $X = \{x_i, x_{i+1}\}_{i \in [k-1]}$. Then it is possible to construct a set equal to $X'$ by choosing one element from each set in $X'$. 

Lemma 14. Let \( T, T' \) be single-labelled trees such that \( T \leq T' \), and let \( u, x, y \) be leaf nodes in \( T \) (and hence also in \( T' \)). If \( \text{lca}_T (u, x) < \text{lca}_T (u, y) \) then \( \text{lca}_{T'} (u, x) < \text{lca}_{T'} (u, y) \).

The following two lemmas form the core of our main result.

Lemma 15. If \((X, C)\) is an instance of \( X3C3 \) that allows an exact 3-cover \( C' \) then the corresponding instance \((\{T_1, T_2\}, Y)\) of \textsc{MulsetPCOMP} admits a perfect pruning giving a compatible set of trees.

Proof. Let \( I \subset [k] \) denote the set of \( m \) indices \( i \) of those \( S_i \) we did not choose as part of our exact 3-cover, that is the sets \( S_i \in C - C' \). Let \( \psi : I \rightarrow [m] \) be an arbitrary isomorphism. Prune the tree \( T_1 \) as follows:

- For each \( S_i \) (or equivalently, each subtree rooted at \( \pi_i \)):
  - If \( S_i \subset C' \) then keep the labels of \( S_i \) as the children of \( \pi'_i \) but prune away all other leaf labels in the subtree \( T'_1 \). After pruning there are three leaves (with labels corresponding to the elements of \( S_i \)) as children of \( \pi'_i \), which is itself a child of \( \mu_i \).
  - If \( S_i \not\subset C' \) then prune away all leaf labels in \( T'_1 \) except \( a^{\psi(i)}_i \) and \( b^{\psi(i)}_i \). After pruning, the remaining leaves will be children of \( \pi_i \).
- In each \( T'_1 \), prune away \( a^j_i \) if the leaf label appears in \( T'_{1j} \). Since \( m \) sets of \( C \) are not in \( C' \), there are \( m \) leaf labels of the form \( a^j_i \) appearing in \( T'_{1j} \), specifically \( a^\psi(i) \) for the \( m \) values of \( i \) for which \( \psi(i) \) is defined, or equivalently for all \( \psi(i) = 1, \ldots, m \). Hence we must prune away one leaf label in each \( T'_{1j} \).
- Repeat the previous step for each \( T'_{1j} \) - the argument is identical.

Denote the pruned version of \( T_1 \) by \( T'_1 \). Every leaf label of \( X \) appears once; this follows directly from \( C' \) being an exact 3-cover. All other leaf labels appear only once by inspection; hence this is a perfect pruning.

We now prune \( T_2 \) as follows:

- For each \( T'_2 \), prune away all leaf labels except \( a^{\psi^{-1}(j)}_i \) and \( b^{\psi^{-1}(j)}_i \); after pruning \( \rho^j \) has two children, both leaves.
- For each \( T'_{2j} \), we wish to prune this subtree to create a \((k - 1)\) leaf star with all labels \( a^j_i \) for our fixed \( j \) except \( a^{\psi^{-1}(j)}_i \). This is possible due to Lemma 13; from each pair of leaf labels rooted by a child of \( \alpha^j \) we pick one leaf to keep and one to prune away such that we keep one copy of everything except \( a^{\psi^{-1}(j)}_i \).
- Repeat the previous step for each \( T'_{2j} \) - as before the argument is identical.

There is one copy of each leaf label of \( X \) in \( T_2 \), which we do not prune. We prune the leaf labels of \( A \cup B \) in \( T_2 \) so that the leaf labels of \( A \cup B \) in \( T_2' \) are exactly those that do not appear in \( T_2'' \). Hence this is a perfect pruning, which we denote \( T'_2 \).

We now show that \( T'_2 \) can be constructed from \( T'_1 \) by repeated non-leaf edge contractions, which will show \( \{T'_1, T'_2\} \) form a compatible set (with \( T^* := T'_1 \)).

In \( T'_1 \) contract every remaining \( \pi'_i \) (one for each of the \( q \) sets \( S_i \in C' \)) into its parent \( \mu_i \). Furthermore in each \( T_1'' \) and \( T_1'' \), contract every non-leaf edge to create a star rooted at \( \alpha_i \) or \( \beta_i \). By inspection, we can see \( T'_2 \leq T'_1 \), giving our required compatible set of trees. ▶

See Figures 8, 9, 10, and 11 for an example of a pruning as in Lemma 15.
Figure 8 As an illustrated example, a possible perfect pruning $T'_1$ as in Lemma 15. In this example, $S_1, S_k \in C'$, but $S_2 \notin C'$.

Figure 9 A subtree of $T'_1$ in the pruning from Figure 8. In this example, $\psi(1 + \frac{k}{2}) = i$.

Figure 10 A possible perfect pruning $T'_2$, corresponding to the perfect pruning of Figure 8.

Figure 11 A subtree of $T'_2$ in the pruning from Figure 10. Again, here $\psi^{-1}(i) = \frac{i}{2} + 1$.

Lemma 16. If $(X, C)$ is an instance of X3C3 that does not allow an exact 3-cover $C'$ then the corresponding instance $(\{T_1, T_2\}, Y)$ of MULSETPComp does not admit a perfect pruning giving a compatible set of trees.

We omit the proof of Lemma 16 due to space concerns. We now prove Theorem 11.
We define a function \(\text{MUL-trees} T\) associating each that for the MAST problem \([13, 22, 32]\). For all \(\bigcup_{i=1}^{k} D(T_i) = \mathcal{X}\), where \(\mathcal{X} = \bigcup_{i=1}^{k} \mathcal{X}(T_i)\).

We adapt a technique using dynamic programming over \(k\)-tuples of nodes previously used for the MAST problem \([13, 22, 32]\). For all \(k\)-tuples of nodes \((a_1, \ldots, a_k) \in \prod_{i=1}^{k} V(T_i)\), let \(S(a_1, \ldots, a_k) = \bigcup_{i=1}^{k} D(T_i^{a_i})\) denote the set of unique leaf labels that occur in the subtrees \(T_i^{a_i}, \ldots, T_k^{a_k}\), rooted at \(a_1, \ldots, a_k\), respectively.

We aim to find a binary tree \(T\) that is leaf-labelled by \(S(a_1, \ldots, a_k)\) such that each of \(T_i^{a_i}\) for \(i = 1, \ldots, k\) displays \(T_i^{a_i}\). Lemma 18, below, shows that the necessary condition of the existence of such a tree is that \(S(a_1, \ldots, a_k) \cap \mathcal{X}(T_i) \subseteq \mathcal{X}(T_i^{a_i})\) for \(i = 1, \ldots, k\).

**Lemma 18.** Let \(T\) be a binary tree leaf-labelled by \(S(a_1, \ldots, a_k)\). If \(S(a_1, \ldots, a_k) \cap \mathcal{X}(T_i) \not\subseteq \mathcal{X}(T_i^{a_i})\) for some \(i\), \(T_i^{a_i}\) does not display \(T_i^{a_i}\).

**Proof.** Suppose that \(S(a_1, \ldots, a_k) \cap \mathcal{X}(T_i) \not\subseteq \mathcal{X}(T_i^{a_i})\), then there exists \(x \in S(a_1, \ldots, a_k)\) such that \(x \not\in \mathcal{X}(T_i^{a_i})\). Since \(x \in T_i^{a_i}\), we have that \(T_i^{a_i}\) cannot display \(T_i^{a_i}\).

Next, for each \(a_i \in V(T_i)\), let \(P(a_i) = \{\epsilon, a_i, a_i', a_i''\}\), for each \(i = 1, \ldots, k\), where \(a_i'\) and \(a_i''\) denotes the two (unordered) children of vertex \(a_i\), and with \(a_i' = \epsilon\) and \(a_i'' = a_i\) in the case that \(a_i \in L\) is a leaf vertex w.l.o.g. Let \(c_i : P(a_i) \to P(x_i)\) be the involution given by

\[
c_i(\epsilon) = a_i, \quad c_i(a_i) = \epsilon, \quad c_i(a_i') = a_i', \quad \text{and} \quad c_i(a_i'') = a_i'',
\]

associating each \(x \in P(a_i)\) with a complement. Let \(\Pi(a_1, \ldots, a_k) = \prod_{i=1}^{k} P(a_i) - \{\epsilon, \ldots, \epsilon\}, \{a_1, \ldots, a_k\}\), and note that \(|\Pi(a_1, \ldots, a_k)| \leq 4^k\), for all \(a_i \in V(T_i)\), for \(i = 1, \ldots, k\).

We define a function \(W : \prod_{i=1}^{k} V(T_i) \to \{\text{true}, \text{false}\}\) recursively, as follows:

- If \(|S(a_1, \ldots, a_k)| \leq 1\), \(W(a_1, \ldots, a_k) = \text{true}\).
- If \(S(a_1, \ldots, a_k) \cap \mathcal{X}(T_i) \not\subseteq \mathcal{X}(T_i^{a_i})\) for some \(i\), \(W(a_1, \ldots, a_k) = \text{false}\).
- Otherwise,

\[
W(\bar{a}) = \bigvee_{x \in \Pi(\bar{a})} \left( W(x \land W(c(\bar{x})) \land Q(\bar{a}, \bar{x}) \land Q(\bar{a}, c(\bar{x}))) \right)
\]
where $\vec{x} = (x_1, \ldots, x_k)$, $\bar{x}(\vec{x}) = (c_1(x_1), \ldots, c_k(x_k))$ and

$$Q(\vec{x}, \vec{\alpha}) = \bigwedge_{i \neq j} \left( x_i = a_i \land x_j = \epsilon \Rightarrow L(a_i) \cap L(a_j) = \emptyset \lor L(a_i) \cap L(a_j) = \emptyset \right)$$

We may define a partial ordering $\prec$ on $\prod_{i=1}^k (V(T_i) \cup \{\epsilon\})$ by taking

$$(a_1, \ldots, a_k) \prec (a'_1, \ldots, a'_k) \iff a_i = \epsilon \lor a_i \prec a'_i, \quad \text{for all } i = 1, \ldots, k$$

where $\prec$ is the successor relation in $T_i$, with the unique $\prec$-minimum element $(\epsilon, \ldots, \epsilon)$. An example computation of the function $W$ is described in Figure 12.

\begin{figure}[h]
\centering
\begin{tikzpicture}
\node at (0,0) (a) {$2 = \text{val}_1$};
\node at (1,0) (b) {$3 = \text{val}_1$};
\node at (2,0) (c) {$4 = \text{val}_1$};
\node at (3,0) (d) {$5 = \text{val}_1$};
\node at (4,0) (e) {$6 = \text{val}_1$};
\node at (5,0) (f) {$7 = \text{val}_1$};
\node at (6,0) (g) {$8 = \text{val}_1$};
\node at (1,-1) (h) {$a$};
\node at (2,-1) (i) {$a$};
\node at (3,-1) (j) {$b$};
\node at (4,-1) (k) {$b$};
\node at (2,-2) (l) {$T_1$};
\node at (4,-2) (m) {$T_2$};
\path[->] (a) edge (h);
\path[->] (b) edge (i);
\path[->] (c) edge (j);
\path[->] (d) edge (k);
\path[->] (e) edge (l);
\path[->] (f) edge (m);
\path[->] (g) edge (m);
\end{tikzpicture}
\caption{Figure 12 Given the trees $T_1$ and $T_2$, $W(1,6) = \text{true}$ since $W(2,7) \land W(3,8) \leftarrow W(2,7) \land W(4,5) \land W(5,8) = \text{true}$. Note that $W(1,6)$ would also compute $W(2,8) \land W(3,7), W(2,6) \land W(3,6), W(1,7) \land W(6,8)$, and $W(1,8) \land W(6,7)$.}
\end{figure}

> Lemma 19. Let $T_1, T_2, \ldots, T_k$ be a collection of $k$ binary MUL-trees such that $\bigcup_i \mathcal{X}(T_i) = \bigcup_i D(T_i)$. Then $W(a_1, \ldots, a_k) = \text{true}$ if and only if there exists a single-labelled tree $T$ leaf-labelled by $S(a_1, \ldots, a_k)$ such that $T_i^{a_i}$ displays $T|_{\mathcal{X}(T_i)}$, under a mapping that maps $r_i(T) \mapsto x_i$, for all $i = 1, \ldots, k$.

**Proof.** We prove this by induction on $k$. For the base case, suppose that each of $(a_1, \ldots, a_k) = (\epsilon, \ldots, \epsilon)$ is the $\prec$-minimum, so that $T_i^{a_i} = T_i = T_\emptyset$ is the empty tree, for $i = 1, \ldots, k$. In which case $S(\epsilon, \ldots, \epsilon) = \emptyset$, and hence $W(a_1, \ldots, a_k) = \text{true}$ by definition, while (trivially) the empty tree $T = T_\emptyset$ is such that $T_i^{a_i} = T_\emptyset$ displays $T_\emptyset|_{\mathcal{X}(T_i)} = T_\emptyset$. Next, suppose that the result holds for all $(a_1, \ldots, a_k) \prec (a'_1, \ldots, a'_k)$ for some tuple $(a_1, \ldots, a_k)$. We claim that the result holds too for $(a_1, \ldots, a_k)$.

($\Rightarrow$) Suppose that $T$ is as described. Then for each $i = 1, \ldots, k$ there is some subtree $S_i \subseteq T_i^{a_i}$ and some label-preserving isomorphism $f_i : V(S_i) \rightarrow V(T_i|_{\mathcal{X}(T_i)})$.

Let $T'$ and $T''$ denote the left and right subtrees attached at $r(T)$. As $T$ is a single-labelled tree, it follows that $L(T') \cap L(T'') = \emptyset$, as each label occurs exactly once. We can partition each $V(S_i)$ into three parts $L_i = \{v \in V(S_i) : f_i(v) \in T'\}$ and $R_i = \{v \in V(S_i) : f_i(v) \in T''\}$ and $C_i = \{v \in V(S_i) : f_i(v) = r(T)\}$. Note that, since $f_i$ is an isomorphism, if $u \in X$ and $u \prec v$ then $v \in X$, for $X \in \{L_i, R_i\}$.

We have three cases depending on which of these three sets lies the root node $a_i$ of the subtree $T_i^{a_i}$:

- If $a_i \in C_i$, then it follows that either $a_i^l \in L_i$ and $a_i^r \in R_i$, or $a_i^l \in L_i$ and $a_i^r \in R_i$.
- If $a_i^l \in L_i$ and $a_i^r \in R_i$, then $T_i^{a_i^l}$ displays $T_i^{a_i^l}|_{\mathcal{X}(a_i)}$ and $T_i^{a_i^r}$ displays $T_i^{a_i^r}|_{\mathcal{X}(a_i)}$.
- If $a_i^l \in L_i$ and $a_i^r \in R_i$, then $T_i^{a_i^l}$ displays $T_i^{a_i^l}|_{\mathcal{X}(a_i)}$ and $T_i^{a_i^r}$ displays $T_i^{a_i^r}|_{\mathcal{X}(a_i)}$. 

In each case, there is some \( x_i \in \{a^*_i, a^*_k\} \) such that \( T^x_i \) displays \( T^x_i \mid X(a_i) \) and \( T^{c(x_i)}_i \) displays \( T^{c(x_i)}_i \mid X(a_i) \).

- If \( a_i \in L_i \) then it follows that \( L_i = V(S_i) \) and \( C_i = R_i = \emptyset \). Hence we have that \( T^x_i \) displays \( T^x_i \mid X(a_i) \), while \( T^{c(x_i)}_i = T_\emptyset \) (trivially) displays the empty tree \( T^x_i \mid X(a_i) = T_\emptyset \).

- Symmetrically, if \( a_i \in R_i \) then \( T^{c(x_i)}_i \) displays \( T^{c(x_i)}_i \mid X(a_i) \) and \( T^x_i \) displays \( T^x_i \mid X(a_i) \).

In all cases there is some \( x_i \in P(a_i) \) such that \( T^x_i \) displays \( T^x_i \mid X(a_i) \) and \( T^{c(x_i)}_i \) displays \( T^{c(x_i)}_i \mid X(a_i) \), for all \( i = 1, \ldots, k \). Since both \( (x_1, \ldots, x_k), (c_1(x_1), \ldots, c_k(x_k)) \subseteq (a_1, \ldots, a_k) \), it follows from the induction hypothesis that \( W(\vec{x}) \wedge W(\vec{c}(\vec{x})) = \text{true} \).

For all \( i \neq j \), if \( x_i = a_i \) and \( x_j = \epsilon \) then by definition \( a_i \in L_i \subseteq L(T_i^x) \) while \( a_j \in C_j \). If \( a^*_j \in R_j \) then \( L(a_j) \cap L(a^*_j) = \emptyset \), otherwise \( a^*_j \in R_j \) and so \( L(a_j) \cap L(a^*_j) = \emptyset \), since \( R_i \subseteq L(T_i) \) and \( L(T_i) \cap L(T_j) = \emptyset \). If \( x_i = a_i \) and \( x_j = a^*_j \) then \( a_i \in L_i \subseteq L(T_i) \) while \( a^*_j \in R_j \) and \( a^*_j \subseteq R_j \subseteq L(T_j) \). From which it follows that \( L(a_j) \cap L(a^*_j) = \emptyset \).

Similarly, if \( x_i = a_i \) and \( x_j = a^*_j \) then it follows \( L(a_j) \cap L(a^*_j) = \emptyset \). This is to say that \( Q(\vec{a}, \vec{x}) = \text{true} \), and by the same argument, so too that \( Q(\vec{a}, \vec{c}(\vec{x})) = \text{true} \).

Hence, by definition, \( W(a_1, \ldots, a_k) = \text{true} \), as required.

\((\Rightarrow)\) Suppose that \( W(a_1, \ldots, a_k) = \text{true} \), then there are two possible cases:

\((i)\) \(|S(a_1, \ldots, a_k)| \leq 1\); \((ii)\) there is some \( (x_1, \ldots, x_k) \in \Pi(a_1, \ldots, a_k) \) such that \( W(\vec{x}) \wedge W(\vec{c}(\vec{x})) \wedge Q(\vec{x}) = \text{true} \).

\((i)\) If \( S(a_1, \ldots, a_k) = \emptyset \) then we may take \( T^x_i \) to display \( T = T_\emptyset \) as the empty tree for all \( i = 1, \ldots, k \). Otherwise, if \( S(a_1, \ldots, a_k) = \{x\} \) is a singleton then we may take \( T \) to be the tree with a single leaf-labelled by \( x \), where it is straightforward to check that \( T^x_i \) can display \( T_i \mid X(a_i) \) for all \( i = 1, \ldots, k \).

\((ii)\) It follows from the induction hypothesis that there exist single-labelled trees \( T_x \) and \( T_y \), leaf-labelled by \( S(x_1, \ldots, x_k) \) and \( S(y_1, \ldots, y_k) \), respectively, such that \( T^x_i \) displays \( T_i \mid X(T_x) \) and \( T^y_i \) under a mapping that maps \( r_i(T_x) \mapsto x_i \), and displays \( T_y \mid X(T_y) \) under a mapping that maps \( r_i(T_y) \mapsto y_i \), for all \( i = 1, \ldots, k \), where \( y_i = c(x_i) \). If \( L(T_x) \cap L(T_y) = \emptyset \) then we construct a new tree \( T \) by connecting the roots \( r(T_x) \) and \( r(T_y) \) of \( T_x \) and \( T_y \) to a common (new) root node \( r \).

Otherwise since \( Q(\vec{a}, \vec{x}) \wedge Q(\vec{a}, \vec{c}(\vec{x})) = \text{true} \) it follows that either \( L(T_x) \cap L(T_y) = \emptyset \) or \( L(T_x) \cap L(T_y) = \emptyset \). In the first case we can construct a new tree \( T \) by merging \( T_x \) with the left sub-tree of \( T_y \), while in the latter case we can construct \( T \) by merging \( T_x \) with the left sub-tree of \( T_y \).

In all cases, we have that \( T \) is a single-labelled tree, as required, as it remains to show that \( T^x_i \) displays \( T_i \mid X(T_x) \) for each \( i = 1, \ldots, k \):

- If \( x_i = a_i \) then \( T^x_i \) displays \( T_i \mid X(T_x) \), mapping \( r(T_x) \mapsto a_i, r(T_y) \mapsto a_i \), and \( r = a_i \).
- If \( x_i = a^*_i \) then \( T^{c(x_i)}_i \) displays \( T_i \mid X(T_y) \), mapping \( r(T_x) \mapsto a^*_i, r(T_y) \mapsto a^*_i \), and \( r = a_i \).
- If \( x_i = a_i \) then \( T^{c(x_i)}_i \) displays \( T_i \mid X(T_y) \) under the mapping that maps \( r(T_x) \mapsto a_i \).
- If \( x_i = \epsilon \) then \( T^{c(x_i)}_i \) displays \( T_i \mid X(T_y) \) under the mapping that maps \( r(T_y) \mapsto a_i \).

Hence, it follows from induction that \( W(a_1, \ldots, a_k) = \text{true} \) if and only if there is some tree \( T \) leaf-labelled by \( S(a_1, \ldots, a_k) \) such that \( T^x_i \) displays \( T_i \mid X(T_x) \) under a mapping that maps \( r_i(T) \mapsto x_i \), for each \( i = 1, \ldots, k \), as required. \(\square\)

Lemma 19 provides us with a criterion for deciding the \textsc{MULSETPC} problem for a given collection of binary MUL-trees, that can be computed in polynomial-time in the size of the trees, for any fixed number of trees, and scales exponentially with the number of trees.
Theorem 20. Let $T_1, T_2, \ldots, T_k$ be a collection of $k$ binary MUL-trees such that $\bigcup_{i} \mathcal{X}(T_i) = \bigcup_{i} D(T_i)$. Then MULSETPC for this instance can be solved in $O(k^2 \cdot 4^k \prod_{i=1}^{k} (|T_i| + 1)) = O(4^k \prod_{i=1}^{k} |T_i|)$ time.

Proof. Based on Lemma 19, it is sufficient to compute $W(r(T_1), \ldots, r(T_k))$, since $T_i^{r(T_i)} = T_i$, by definition. We can compute $W$ via dynamic programming as outlined in Algorithm 1, which will return true if $T_1, \ldots, T_k$ display a single labelled tree.

Algorithm 1 Recursive dynamic programming algorithm for $W(a_1, \ldots, a_k)$.

1: let $S = S(a_1, \ldots, a_k)$
2: if $|S| \leq 1$ then return true
3: else if $S \cap \mathcal{X}(T_i) \not\subseteq \mathcal{X}(T_i^{a_i})$ for some $i = 1, \ldots, k$ then return false
4: else
5: for $(x_1, \ldots, x_k) \in \Pi(a_1, \ldots, a_k)$ do
6: if $W(\vec{x}) \land W(\vec{c}(\vec{x})) \land Q(\vec{a}, \vec{x}) \land Q(\vec{a}, \vec{c}(\vec{x}))$ then return true
7: return false

For the time complexity, we can use memoization to store the values of $W$ in a table with at most $O(|\Pi_{i=1}^{k} (|T_i| + 1))$ entries. Furthermore, we require at most $O(k^2 \cdot 4^k)$ time to compute the value of each entry $W(a_1, \ldots, a_k)$, since $|\Pi(a_1, \ldots, a_k)| = |P(a_1)| \times \cdots \times |P(a_k)| \leq 4^k$, while $Q(\vec{a}, \vec{x})$ and $Q(\vec{a}, \vec{c}(\vec{x}))$ can each be computed in quadratic time. Hence, the running time is $O(k^2 \cdot 4^k \prod_{i=1}^{k} (|T_i| + 1)) = O(4^k \prod_{i=1}^{k} |T_i|)$, as required.

6 Conclusions

The above results resolve an open problem posed in [15, 16] as to whether the MULSETPC problem remains NP-complete when the number of MUL-trees is constant. According to Theorems 9 and 10, two MUL-trees are sufficient for NP-completeness, even with each label appearing at most twice in any tree and either the height or the degree constant. Theorems 11 and 12 extend this result and show that the more general MULSETPCComp problem also remains NP-complete even when the number of MUL-trees is constant. Theorem 9 is tight in the sense that, if we restrict our attention to MUL-trees in which each label appears uniquely in at least one tree, we obtain a polynomial-upper bound for a fixed number of trees (Theorem 20). However, Theorem 12 suggests the algorithm presented in Theorem 20 for MULSETPC cannot be directly generalised to solve equivalent MULSETPCComp instances in polynomial time, unless $P = NP$.

The above results also suggest two new open problems. Firstly, is it possible to improve Theorem 11 to show that MULSETPCComp is still NP-complete when restricted to MUL-trees with multiplicity 2? Secondly, what can be said about the complexity of MULSETPC and MULSETPCComp for instances in which the multiplicity is not restricted but the number of leaf labels that may appear more than once is restricted? That is, for each MUL-tree, $k$ leaf labels may appear an unbounded number of times in the tree, whereas all other labels appear at most once. For which values of $k$ are these subproblems NP-complete? This question is interesting because of its connection to the instances investigated in Section 5.

References


14:18  MUL-Tree Pruning for Consistency and Compatibility


Linear-Time Computation of Cyclic Roots and Cyclic Covers of a String

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Abstract
Cyclic versions of covers and roots of a string are considered in this paper. A prefix V of a string S is a cyclic root of S if S is a concatenation of cyclic rotations of V. A prefix V of S is a cyclic cover of S if the occurrences of the cyclic rotations of V cover all positions of S. We present O(n)-time algorithms computing all cyclic roots (using number-theoretic tools) and all cyclic covers (using tools related to seeds) of a length-n string over an integer alphabet. Our results improve upon O(n log log n) and O(n log n) time complexities of recent algorithms of Grossi et al. (WALCOM 2023) for the respective problems and provide novel approaches to the problems. As a by-product, we obtain an optimal data structure for Internal Circular Pattern Matching queries that generalize Internal Pattern Matching and Cyclic Equivalence queries of Kociumaka et al. (SODA 2015).

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1 Introduction

Cyclic strings have many real-world applications, such as in bioinformatics [2, 3, 17, 19] and image processing [1, 27, 28, 29]. In particular, they are used for detecting DNA viruses with circular structures [30, 31]. In particular, cyclic strings were studied in the context of circular pattern matching [10, 14, 23, 24].

In this paper, we investigate the complexity of two problems related to cyclic strings. The first one is a cyclic variant of the problem of computing the roots of a string S, i.e., strings U such that S = U^k for some integer k. The second one is a cyclic variant of the problem of computing the covers of a string S, i.e., strings C whose occurrences cover the...
whole string $S$. The standard roots of a string can be easily computed in linear time using a folklore algorithm. Moore and Smyth [25, 26] gave a linear-time algorithm computing all standard covers of a string. However, the cyclic versions of these problems are more difficult.

We say that a string $V$ is a \( \text{cyclic rotation} \) of a string $U$ if there exist strings $X$ and $Y$ such that $U = XY$ and $V = YX$. A string $U$ has a \text{circular occurrence} in a string $T$ at position $i$ if a rotation of $U$ has a (standard) occurrence in $T$ at position $i$.\(^1\) By $\text{CircOcc}(U, T)$ we denote the set of circular occurrences of $U$ in $T$. Moreover, we denote

\[
\text{Covered}(U, T) = \bigcup_{i \in \text{CircOcc}(U, T)} [i..i + |U|).
\]

\textbf{Definition 1.} A string $U$ is a \text{cyclic cover} of a string $S$ if $\text{Covered}(U, S) = [0..|S|)$. A string $U$ is a \text{cyclic root} of a string $S$ if $S = U_1 \cdots U_k$, where each $U_i$ is a cyclic shift of $U$.

Note that if $U$ is a cyclic root (cyclic cover) of $S$, then the prefix $S[0..|U|]$ is also a cyclic root (cyclic cover, respectively) of $S$.

\textbf{Example 2.} The lengths of the cyclic roots of the Thue–Morse word $\text{abbaabbaababa}$ are $2, 4, 8, 16$.

\textbf{Example 3.} The string $\text{ab}$ is a cyclic cover of each Fibonacci string of length at least 2, e.g., of the string $\text{Fib}_5 = \text{abababa}$. (See [13] for a definition of Fibonacci strings and their properties.) However, it is not a cyclic root of any Fibonacci string longer than 2. Another example of a cyclic cover of a string is illustrated in Figure 1.

We consider the following problems.

\begin{table}[h]
\begin{tabular}{|l|}
\hline
\textbf{CyclicRoots} \\
\textbf{Input:} A string $S$ of length $n$. \\
\textbf{Output:} The lengths of all cyclic roots of $S$. \\
\hline
\end{tabular}
\end{table}

\begin{table}[h]
\begin{tabular}{|l|}
\hline
\textbf{CyclicCovers} \\
\textbf{Input:} A string $S$ of length $n$. \\
\textbf{Output:} The lengths of all cyclic covers of $S$. \\
\hline
\end{tabular}
\end{table}

\textbf{Our results}

We show linear-time algorithms for both problems. We assume the word-RAM model of computation and that the string $S$ is over an integer alphabet $\{0, \ldots, \Theta(n)\}$.

\textsuperscript{1} We assume that the positions of a string $T$ are numbered 0 through $|T| - 1$, where $|T|$ is the length of $T$. 
Previous results

Recently, Grossi et al. [16] presented an $O(n \log \log n)$-time algorithm for CyclicRoots (named cyclic factorization there) and an $O(n \log n)$-time algorithm for CyclicCovers.

Remark 4. A completely different problem of covering a cyclic string with a standard string cover was considered in [11, 12]. Another different problem also known under the name “cyclic covers”, related to the shortest superstring problem, was considered in [4, 5, 6].

Our approach

In the case of cyclic covers, we use a recursive algorithm whose general structure partially resembles the structure of the linear-time algorithm for computing seeds from [21]. For this, we need to explore combinatorics of circular occurrences, which is different from that of standard occurrences. The “working horse” of the algorithm (non-recursive parts) is the computation of long cyclic covers, which is based on an efficient implementation of internal circular pattern matching queries. Such queries require to find all circular occurrences of one substring of a text in another substring: the set of occurrences can be represented compactly if the ratio of lengths of the two strings is constant. An auxiliary contribution of our paper is an optimal implementation of these queries (constant-time after linear-time preprocessing).

Also in the case of cyclic roots we use internal queries on strings. Our algorithm is based on number-theoretic tools and fast internal queries for cyclic equivalence, which ask if two substrings of a given string are rotations of each other [20, 22].

Notations

For a string $S$, by $S[0], \ldots, S[|S| - 1]$ we denote its respective letters. By $S[i..j]$ we denote a substring $S[i] \cdots S[j - 1]$; similarly, we define substrings $S[i..j]$, $S(i..j)$ and $S(i..j)$. We say that $p$ is a period of a string $S$ if $S[i] = S[i + p]$ holds for all $i \in [0..|S| - p)$. By $\text{per}(S)$, we denote the smallest period of $S$, called the period of $S$. For a string $S$ and integer $x \in [0..|S|)$, by $\text{rot}_x(S)$ we denote the rotation $S[x..|S|] \cdot S[0..x)$ of $S$ obtained from $S$ by moving the prefix of $S$ of length $x$ to the end.

A length-$m$ string $P$ has an occurrence in a string $T$ at position $i$ if $T[i..i + m) = P$. By $\text{Occ}(P,T)$ we denote the set of starting positions of occurrences of $P$ in $T$.

2 Internal Circular Pattern Matching and CyclicCovers in $O(n \log n)$ Time

We introduce the following generalization of Internal Pattern Matching queries from [22].

**Simple Internal Circular Pattern Matching Queries (Simple InternalCPM)**

**Input:** A string $S$ of length $n$.

**Queries:** Given two substrings $P$ and $T$ of $S$ such that $|T| \leq 2|P|$, report the leftmost and the rightmost circular occurrence of $P$ in $T$.

Remark 5. If we know how to compute the leftmost circular occurrence, then the rightmost circular occurrence can be computed analogously (it suffices to reverse the strings).

The theorem below can be obtained using the methods from [7, 8]. We give its proof in Section 6.

Theorem 6. The Simple InternalCPM queries can be answered in $O(1)$ time after $O(n)$-time preprocessing.
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Remark 7. In Section 6, we obtain a version of the queries in which a constant-sized representation of all circular occurrences is computed in constant time (still if $|T| \leq 2|P|$).

Below, we apply Simple InternalCPM queries to a version of the CyclicCovers problem that is used in our $O(n)$-time algorithm for CyclicCovers. The following lemma generalizes [16, Lemma 13]; in this section, it will be applied in a simpler setting.

Lemma 8. After $O(n)$-time preprocessing of a string $S$ of length $n$, for any substrings $C$ and $W$ of $S$, we can test if $C$ is a cyclic cover of $W$ in $O(|W|/|C|)$ time.

Proof. Let $p = |C|$ and $m = |W|$. Consider substrings $V_i$ equal to $W[ip..(i+2)p]$ for $i \in [0..[m/p]-1]$ and $W[ip..m]$ for $i = [m/p]-1$. For each substring $V_i$, we use a Simple InternalCPM query to compute the leftmost and the rightmost circular occurrence of $C$. Then, we check whether these occurrences, interpreted as occurrences in $W$, collectively cover all positions in $W$. The time complexity is $O(m/p)$ after the preprocessing of Theorem 6.

Lemma 8 implies a simple $O(n \log n)$-time algorithm for the CyclicCovers problem.

Corollary 9. The CyclicCovers problem can be solved in $O(n \log n)$ time.

Proof. We apply Lemma 8 for $W = S$ iterating with $C$ over all non-empty prefixes of $S$. The total time complexity is $O(n + \sum_{i=1}^{n} \frac{n}{i}) = O(n \log n)$.

3 Quasi-Covers

We reduce our problem to the computation of the substrings called quasi-covers; see Figure 2.

Definition 10. A string $V$ is a quasi-cover of a substring $W$ of the string $S$ if $V$ is a prefix of $S$ and a cyclic cover of a substring $Y = W[i..j]$ such that $i < |V|$ and $j > |W| - |V|$.

Figure 2 Example of a quasi-cover $aab$ of the substring $W$ of $S$. By definition, $V$ is a prefix of the whole string $S$. Observe that $aab$ is not a cyclic cover of $W$.

Henceforth, we fix the string $S$ and consider quasi-covers of its substrings. Let $W$ be a substring and $I$ be an interval, $I \subseteq [1..|W|]$. We denote by $Q$-COVERS$_I(W)$ the set of all lengths of quasi-covers of substring $W$ with lengths in $I$. Furthermore, for $k \in [1..|W|]$ we denote

$$Q$-COVERS(W) = Q$-COVERS_{[1..|W|]}(W),$$

$$Q$-COVERS_{\leq k}(W) = Q$-COVERS_{[1..k]}(W),$$

$$Q$-COVERS_{> k}(W) = Q$-COVERS_{[k..|W|]}(W),$$

$$Q$-COVERS_{=}k(W) = Q$-COVERS_{[k..k]}(W).$$

A prefix of $S$ is its cyclic cover if and only if it is a quasi-cover of $S$ and a rotation of a suffix of $S$. We use the following queries (generalized by Simple InternalCPM queries):
Cyclic Equivalence Queries (CycEq)
Input: A string \( S \) of length \( n \).
Queries: Given two substrings \( U \) and \( V \) of \( S \), check if \( V \) is a rotation of \( U \).

\[ \text{Theorem 11 ([20, 22])}. \] The CycEq queries can be answered in \( O(1) \) time after \( O(n) \)-time preprocessing.

Using CycEq queries, we can compute in \( O(n) \) time all prefixes which are rotations of the corresponding suffixes of the string. This yields the following observation.

\[ \text{Observation 12}. \] The CyclicCovers problem for a string \( S \) reduces in linear time to the computation of Q-Covers(S).

We will later show how the set Q-Covers(S) can be computed based on recursive calls to Q-Covers(W) for substrings \( W \) of \( S \).

3.1 Quasi-Covers and Substring Complexity
The substring complexity of a length-\( m \) string \( W \) is a function that maps each length \( k \in [1 \ldots m] \) to the number \( |\text{SUB}_k(W)| \) of distinct length-\( k \) substrings of \( W \). We further define \( \beta_k(W) = |\text{SUB}_k(W)| + k - 1 \). The term \( k - 1 \) is added because the sequence \( (|\text{SUB}_k(W)|)_{k=1}^m \) does not need to be monotone in general; the resulting sequence \( (\beta_k(W))_{k=1}^m \) is now non-decreasing and its monotonicity will be useful later. For a string family \( S \), let us denote by \( \|S\| \) the sum of lengths of strings in \( S \).

\[ \text{Observation 13}. \] Let \( V \) be a quasi-cover of a substring \( W \) of \( S \). If a substring \( W' \) of \( W \) satisfies \( |W'| \geq 2|V| - 1 \), then \( V \) is a quasi-cover of \( W' \).

\[ \text{Lemma 14}. \] Given a length-\( m \) substring \( W \) and an integer \( k \in [1 \ldots m] \), we can compute in \( O(m) \) time a family \( G_k(W) \) of substrings of \( W \) such that \( \|G_k(W)\| \leq \beta_k(W) \) and \( \text{Q-Covers} \leq [k/4] \)(W) = \( \bigcap_{W' \in G_k(W)} \text{Q-Covers} \leq [k/4] \)(W').

Proof. It was shown in [21] that one can construct in linear time a string family, denoted in [21] as \( \text{COMPR}_t \), such that \( \|\text{COMPR}_t\| \leq \beta_{2t-1}(W) \) and the strings in \( \text{COMPR}_t \) contain all length-\( t \) substrings of \( W \). First, we reformulate the corresponding fact from [21] taking \( G_k(W) = \text{COMPR}_{[k/2]} \).

\[ \text{Claim 15 ([21, Lemma 5.4, proof of Lemma 5.3 and proof of Theorem 9 (“Computing S’”))].} \]
Given \( k \in [1 \ldots m] \), we can compute in \( O(m) \) time a string family \( G_k(W) \) such that

\[ ||G_k(W)|| \leq \beta_k(W) \text{ and } \text{SUB}_{[k/2]}(W) = \bigcup_{W' \in G_k(W)} \text{SUB}_{[k/2]}(W'). \]

The next claim turns out to be similar to [21, Lemma 2.2].

\[ \text{Claim 16}. \] If \( t \in [1 \ldots m] \), then

\[ \text{Q-Covers} \leq [t/2] \)(W) = \( \bigcap_{W' \in \text{SUB}_t(W)} \text{Q-Covers} \leq [t/2] \)(W'). \]
Proof. We prove two inclusions separately.

(⊆) If \( V \) is a quasi-cover of \( W \) of length at most \([t/2]\) and \( W' \in \text{SUB}_t(W) \), then Observation 13 implies that \( V \) is a quasi-cover of \( W' \).

(⊇) Assume that \( V \) is a quasi-cover of all \( W' \in \text{SUB}_t(W) \) and \( |V| = \ell \leq [t/2] \). Consider a position \( i \in [\ell - 1 \ldots m - \ell] \) and a substring \( W' = W(i - \ell \ldots i + \ell) \). Note that \( V \) is a quasi-cover of \( W' \) (this follows from Observation 13 because \( W' \) is a substring of some length-\( t \) substring of \( W \)). Thus, there is a circular occurrence of \( V \) in \( W \) that covers the middle position of \( W' \).

Interpreted as a circular occurrence of \( V \) in \( W' \), it covers position \( i \) of \( W' \).

The thesis follows directly from the two claims above, taking \( t = [k/2] \) in the second claim.

Lemma 17. If a string \( W \) has a quasi-cover \( V \) of length \( |V| \geq 2k \), then

\[
|\text{SUB}_{k+1}(W)| \leq \frac{1}{2}|W| + \frac{3}{2}|V|.
\]

Proof. First, we show the following claim (cf. Figure 3).

Claim 18. If a string \( Y \) has a cyclic cover \( V \) of length \( |V| \geq 2k \), then \(|\text{SUB}_{k+1}(V)| \leq \frac{1}{2}(|Y| + |V|)\).

Proof. We denote by \( \text{CSUB}_{k+1}(V) \) the set of distinct length-\((k+1)\) substrings of all rotations of \( V \); note that \(|\text{CSUB}_{k+1}(V)| \leq |V|\).

For each \( i \in \text{Circ Occ}(V,Y) \), let us mark positions \( j \in [i \ldots i + |V| - k) \); observe that if a position \( j \) is marked, then \( Y[j \ldots j + k] \in \text{CSUB}_{k+1}(V) \).

Let \( Y' \) be the prefix of \( Y \) of length \(|Y| - |V|\). We partition \( Y' \) into inclusion-wise maximal intervals of marked positions and inclusion-wise maximal intervals of unmarked positions.

Each interval \( I_2 \) of unmarked positions is preceded by an interval \( I_1 \) of marked positions, where \(|I_1| \geq |V| - k \geq k \geq |I_2|\) (otherwise, \( V \) would not be a cyclic cover of \( Y \)).

Hence, at most half of positions of \( Y' \) are unmarked, which is \(|Y| - |V|)/2 \). Each length-\((k+1)\) substring starting at marked position belongs to \( \text{CSUB}_{k+1}(V) \). Hence, \(|\text{SUB}_{k+1}(V)| \leq (|Y| - |V|)/2 + |\text{CSUB}_{k+1}(V)| \leq (|Y| + |V|)/2 \).

If \( V \) is a quasi-cover of \( W \) then \( V \) is a cyclic cover of \( Y = W[i \ldots j] \) with \( i < |V| \) and \( j > |V| \). We have, due to inequality \( \frac{1}{2}(|W| - |Y|) < |V| \),

\[
|\text{SUB}_{k+1}(W)| \leq |\text{SUB}_{k+1}(Y)| + |W| - |Y| < |\text{SUB}_{k+1}(Y)| + \frac{1}{2}(|W| - |Y|) + |V|.
\]

Now, Claim 18 implies \(|\text{SUB}_{k+1}(Y)| \leq \frac{1}{2}(|Y| + |V|)\) and thus \(|\text{SUB}_{k+1}(W)| \leq \frac{1}{2}|W| + \frac{3}{2}|V|\).
Example 19. Let \( \{a_1, a_2, \ldots, a_{2k-1}\}, \{b_1, b_2, \ldots, b_{2k}\}, \{c_1, c_2, \ldots, c_{2k-1}\} \) be disjoint sets, and
\[
X = a_1 a_2 \cdots a_{2k-1}, \quad V_1 = b_1 b_2 \cdots b_k, \quad V_2 = b_{k+1} b_{k+2} \cdots b_{2k}, \quad Y = c_1 c_2 \cdots c_{2k-1}.
\]
Then \( V = V_1 V_2 \) is a quasi-cover of \( W = X V_1 V_2 V_1 Y \), with \(|V| = 2k\) and \(|W| = 8k - 2\). All length-\((k + 1)\) substrings of \( W \) are different. Hence:
\[
|\text{SUB}_{k+1}(W)| = |W| - k = \frac{1}{2}|W| + \frac{3}{2}|V| - o(|W|).
\]

We use the following crucial property of quasi-covers.

Lemma 20 (Work-Reduction Lemma). For a length-\(m\) substring \( W \) and \( k \in [0..m) \), if \( \beta_{k+1}(W) > \frac{5}{6}m \), then
\[
\text{Q-COVERS}_{[2k...(m/6)]}(W) = \emptyset.
\]

Proof. The proof is by contradiction. Suppose that \( V \in \text{Q-COVERS}_{[2k...(m/6)]}(W) \) and \( V \) is a cyclic cover of \( Y = W[i..j] \) with \( i < |V| \) and \( j > m - |V| \). Due to Lemma 17 and inequality \( k \leq |V|/2 \),
\[
|\text{SUB}_{k+1}(W)| + k \leq \frac{1}{2}|W| + \frac{3}{2}|V| + \frac{1}{2}|V| = \frac{3}{2}|W| + 2|V| \leq \frac{3}{2}m + 2 \cdot \frac{m}{6} = \frac{5}{6}m.
\]
This contradicts our assumption that \( \beta_{k+1}(W) = |\text{SUB}_{k+1}(W)| + k > \frac{5}{6}m \).  ▶

4 Solution to CyclicCovers Problem

Our algorithm is recursive; its non-recursive parts correspond to (simple) fast computation of length-limited cyclic covers. We say that an interval \( I = [a..b] \) of positive integers is balanced if \( b = O(a) \).

Lemma 21. After \( O(n) \)-time preprocessing, for a balanced interval \( I = [a..b] \) and a length-\(m\) substring \( W \), the set \( \text{Q-COVERS}_I(W) \) can be computed in \( O(m) \) time.

Proof. We consider each length \( \ell \in I \) separately. Let \( C = S[0..\ell] \). We use two \textsc{Simple InternalCPM} queries to check if \( C \) has a circular occurrence starting within the first \( \ell \) positions of \( W \) and a circular occurrence ending within the last \( \ell \) positions of \( W \). If any of these two conditions does not hold, \( C \) is not a quasi-cover of \( W \). Otherwise, we use Lemma 8 to check if \( C \) is a cyclic cover of the substring of \( W \) spanned by the first and the last circular occurrence of \( C \) in \( W \) that were discovered in the previous step. The total time complexity is \( O(\sum_{i \in I} \frac{m}{a}) = O(\sum_{i \in I} \frac{m}{b/a}) = O(m \cdot |I|/a) = O(m \cdot b/a) = O(m) \), after \( O(n) \)-time preprocessing in Theorem 6 and Lemma 8.  ▶

Our solution is based on Lemmas 14, 20, and 21. We use a recursive approach that was initially developed for seeds computation; see [21].

Theorem 22. The CyclicCovers problem can be solved in \( O(n) \) time.

Proof. We run the recursive function \textsc{ComputeQuasiCovers} (Algorithm 1) initially for \( W = S \).

Correctness. In the base case, where \( \beta_1(W) > \frac{5}{6}m \), there are more than \( \frac{5}{6}m \) different letters in \( W \), and then Lemma 20 implies \( \text{Q-COVERS}_{[m/6]}(W) = \emptyset \).

In the recursive step, we reduce the computation of quasi-covers to the ones with lengths in two balanced intervals, \( J_1 = ([k/4]..2k) \) and \( J_2 = ([m/6]..m] \), and the ones (the set \( Q \)) with sufficiently small lengths (at most \([k/4]\)). By Lemmas 14 and 20, the algorithm returns precisely the set \( \text{Q-COVERS}(W) \).
Linear-Time Computation of Cyclic Roots and Cyclic Covers of a String

Algorithm 1 ComputeQuasiCovers(W).

Input: A substring W of length m.
Output: The set Q-Covers(W) of lengths of quasi-covers of W.

Compute $\beta_i(W)$ for all $i \in [1..m]$ if $\beta_i(W) > \frac{5}{6} m$ then return $\text{Q-Covers}_{\lfloor(m/6)\rfloor..m}(W)$ \hspace{1cm} \text{see Lemma 20}

$k := \max \{i \in [1..m] : \beta_i(W) \leq \frac{5}{6} m\}$

Let $G_k(W)$ be the set of fragments as in Lemma 14

foreach string $W' \in G_k(W)$ do
\hspace{1cm} $Q_{W'} := \text{ComputeQuasiCovers}(W')$ \hspace{1cm} \text{Recursive call}
\hspace{1cm} $Q := \bigcap_{W' \in G_k(W)} Q_{W'} \cap [1..\lceil k/4 \rceil]$ \hspace{1cm} \text{Q-Covers}_{\lfloor k/4 \rfloor}(W) (Lemma 14)
\hspace{1cm} $J_1 := ([\frac{k}{2}]..2k), J_2 := ([\frac{m}{6}]..m]$ \hspace{1cm} \text{J1, J2 are balanced intervals}
\hspace{1cm} return $Q \cup \text{Q-Covers}_{J_1}(W) \cup \text{Q-Covers}_{J_2}(W)$ \hspace{1cm} \text{Q-Covers}_{2k..\lceil m/6 \rceil}(W) = \emptyset$

Complexity. To bound the running time, denote by $T(m)$ the maximum number of operations performed by the algorithm for a substring $W$ of length $m$. The sequence $\beta_i(W)$ for a length-$m$ substring $W$ of $S$ can be computed in $O(m)$ time [21, Lemma 5.1]. Due to Lemmas 14 and 21,

$$T(m) = O(m) + \sum_i T(m_i), \quad \text{where} \quad \sum_i m_i \leq \frac{5}{6} m.$$ 

This recurrence yields $T(m) = O(m)$.

Due to Observation 12, the CYCLICCOVERS problem can be reduced in linear time to the computation of all quasi-covers. Finally, Lemma 21 requires $O(n)$-time preprocessing of $S$. This completes the proof.

5 Solution to CyclicRoots Problem

We denote by $\sigma_0(n) = \sum_{p|n} 1$ the number of divisors of $n$ and by $\sigma_1(n) = \sum_{p|n} p$ the sum of divisors of $n$. We use the following known estimations: $\sigma_0(n) = 2^{O(\log n / \log \log n)}$ [18, §18.1] and $\sigma_1(n) = O(n \log \log n)$ [18, §22.9]. They directly imply the following fact.

Fact 23.

- $\sigma_0(n) = o(\sqrt{n} / \log n)$ and $\log \sigma_0(n) = O(\log n / \log \log n)$
- $\sigma_1(n) = \sum_{p|n} \frac{n}{p} = O(n \log \log n)$

Using CYCLEQ queries (Theorem 11), we derive the following subroutine:

Observation 24. After linear-time preprocessing of a string $S$, we can test if $S[0..p]$ is a cyclic root of a substring $W$ of $S$ in $O(|W|/p)$ time.

In particular, in [16] the CYCLICROOTS problem was solved in $O(\sigma_1(n)) = O(n \log \log n)$ time (cf. Fact 23) by using $\frac{n}{p}$ CYCLEQ queries for each divisor $p$ of $n$.

Let us now develop an $O(n)$-time solution. We reduce testing if $S[0..p]$ is a cyclic root of the whole text to testing if $S[0..p]$ is a cyclic root of each substring $F$ in a suitably chosen family $\mathcal{F}$ of substrings.
The intuition behind this improvement is as follows. It turns out that the asymptotic upper bound on \( \sigma_1(n) \) significantly depends on a few largest divisors. In the \( O(n \log \log n) \)-time algorithm, this corresponds to the smallest lengths \( p \) of the candidate cyclic root. Hence, for small \( p \), we will adopt a different approach.

The factorization of \( S \) into length-\( q \) substrings, for \( q \mid n \), will be called the \( q \)-factorization.

\[ \text{Observation 25. If } S \text{ has a cyclic root of length } p, \text{ then its } (k \cdot p)\text{-factorization } F \text{ contains at most } p^k \text{ distinct substrings, consequently the number of distinct factors in } F \text{ is at most } \min\left(p^k, \frac{n}{k \cdot p}\right). \]

Thus, if the number of different factors in the \((k \cdot p)\text{-factorization}\) is greater than \( p^k \), then we know that \( S \) does not have a cyclic root of length \( p \).

Otherwise, if \( k \) is small enough, the number of different substrings in the \((k \cdot p)\text{-factorization}\) will be smaller than \( n/(k \cdot p) \), and we can check each of them using \text{CycEq} queries in \( O(k) \) time. On the other hand, if \( k \) is large enough, then the \( O(n/(k \cdot p)) \) work spent on computing the factorization will be much less than \( O(n/p) \).

\textbf{Algorithm 2 CyclicRoot}(\( S, p \)): Does \( S \) have a cyclic root of length \( p \mid n \)?

\[
k := \max(\frac{1}{\log_p n}, 1)
\]

Let \( S = S_1 \cdot S_2 \), where \( |S_2| = n \mod (k \cdot p) \)

\( F := \text{all distinct factors in the } (k \cdot p)\text{-factorization of } S_1 \quad \text{\( O(n/(k \cdot p)) \) time [15]} \)

if \( |F| > p^k \) then return \text{NO}  
if \( |S_2| > 0 \) then \( F := F \cup \{S_2\} \quad \text{\( |F| = O\left(\min\left(p^k, \frac{n}{k \cdot p}\right)\right)\)} \)

foreach \( F \in F \) do 

if \( S[0 \ldots p) \) is not a cyclic root of \( F \) then return \text{NO} 

return \text{YES} 

\[ \text{Observation 24} \]

\[ \text{Theorem 26. The CyclicRoots problem can be solved in } O(n) \text{ time.} \]

\textbf{Proof.} We use Algorithm 2. After \( O(n) \)-time preprocessing, all different substrings in \( F \) can be found in \( O(n/(k \cdot p)) \) time using deterministic substring hashing [15]. By Observation 24, after \( O(n) \)-time preprocessing, we can test in \( O(k) \) time for each \( F \in F \) if \( S[0 \ldots p) \) is its cyclic root; this sums up to \( O(k \cdot \min(\frac{n}{k \cdot p}, p^k)) \) time. For \( k = \max(\frac{1}{\log_p n}, 1) \), we have

\[
\frac{n}{k \cdot p} = O\left(\frac{n \log p}{p \log n}\right) \quad \text{and} \quad k \cdot \min\left(p^k, \frac{n}{k \cdot p}\right) = O\left(\frac{1}{\log_p n} \cdot \log_p n + \min\left(p, \frac{n}{p}\right)\right) = O(\sqrt{n} \log n).
\]

Thus, after \( O(n) \)-time preprocessing, all calls to the algorithm \text{CyclicRoot}(S, p) for all divisors \( p \) of \( n \) work in total time

\[ O(A(n) + B(n)), \text{ where } A(n) = \sum_{p \mid n} \frac{n \log p}{p \log n} \text{ and } B(n) = \sum_{p \mid n} \sqrt{n} \cdot \log n. \]

\textbf{Estimating } \( B(n) \). \text{ By Fact 23, we have}

\[
B(n) = \sqrt{n} \log n \cdot \sum_{p \mid n} 1 = \sqrt{n} \log n \cdot \sigma_0(n) = o(n).
\]
Estimating $A(n)$. We partition the underlying sum into elements that do not exceed $\sigma_0(n)$ and the remaining elements. The former is bounded from above, due to Fact 23, as:

$$\sum_{p\leq \sigma_0(n)} \frac{n \log p}{p \log n} \leq \sum_{p\leq \sigma_0(n)} \frac{n \log \sigma_0(n)}{p \log n} = \frac{\log \sigma_0(n)}{\log n} \sum_{p\leq \sigma_0(n)} \frac{n}{p} = O \left( \frac{1}{\log \log n} \cdot \sigma_1(n) \right) = O(n).$$

The latter sum is bounded from above by:

$$\sum_{p>\sigma_0(n)} \frac{n \log p}{p \log n} \leq \sum_{p>n} \frac{n}{\sigma_0(n)} \sum_{p=n}^1 = \frac{n}{\sigma_0(n)} \cdot \sigma_0(n) = n.$$

This concludes the complexity analysis of the algorithm.

6 InternalCPM via PILLAR Model

6.1 PILLAR Model

We use the so-called PILLAR model that was introduced in [9]. In this model, we assume that the following primitive queries can be performed efficiently, where the argument strings are represented as substrings of strings in a given collection $X$:

- **Extract**$(U, \ell, r)$: Retrieve the substring $U[\ell \ldots r]$.
- **LCP**$(U, V)$, **LCP**$_R(U, V)$: Find the length of the longest common prefix/suffix of $U$ and $V$.
- **IPM**$(U, V)$: Assuming that $|V| < 2|U|$, compute the starting positions of all exact occurrences of $U$ in $V$, expressed as an arithmetic sequence. If the sequence has at least three terms, its difference equals $per(U)$.
- **Accessa**$(U, i)$: Retrieve the letter $U[i]$.
- **Length**$(U)$: Compute the length $|U|$ of the string $U$.

The runtime of algorithms in this model can be expressed in terms of the number of primitive PILLAR operations. The following result combines several known techniques to obtain constant-time implementations of all PILLAR operations in the standard setting, efficient implementations of the PILLAR operations in other settings, including a dynamic and a compressed setting, are also known; cf. [9].

**Theorem 27** ([9, Theorem 7.2]). After an $O(n)$-time preprocessing of a collection of strings of total length $n$ over an integer alphabet, each PILLAR operation can be performed in $O(1)$ time.

6.2 Interval Chains and PairMatch problem

For an integer set $A$ and an integer $r$, let $A \oplus r = \{a + r : a \in A\}$. An interval chain is a set of the form $I \cup (I \oplus q) \cup (I \oplus 2q) \cup \cdots \cup (I \oplus aq)$ for an interval $I$ and non-negative integers $a$ and $q$. In particular, a single interval is an interval chain (with $a = 0$).

First, we introduce an auxiliary operation **PAIRMATCH**. Denote by **PAIRMATCH**$(T, P, i, j)$ the set of all circular occurrences of $P$ in $T$ such that position $i$ in $T$ is aligned with position $j$ in $P$ (see also Figure 4):

**PAIRMATCH**$(T, P, i, j) = \{p \in (i-m..i] : T[p..p+m] = \text{rot}_x(P), i-p = (j-x) \mod m\}$.

In particular **PAIRMATCH**$(T, P, i, 0)$ is the set of circular occurrences of $P$ such that the leftmost position of $P$ is aligned with position $i$ in $T$. 
Figure 4 Let us put original numbers in positions of the pattern $P$; they are moved after rotation of $P$. Assume that $p = 9$ is a position of a circular occurrence of $P$ in $T$ such that $p \in \operatorname{Occ}(\text{rot}_7(P), T)$. Then, in particular, $p \in \operatorname{PairMatch}(T, P, 17, 2)$. In this case, $x = 7$ and $i - p = 8 = (j - x) \mod 13$. We also have $p \in \operatorname{PairMatch}(T, P, 15, 0)$.

The following lemma is a consequence of [7, Lemma 10], where the PILLAR model was not used explicitly. A similar fact was shown in [16, Lemmas 5 and 6]. We include its proof for completeness.

▶ Lemma 28. For any given $i, j$, the set $\operatorname{PairMatch}(T, P, i, j)$, represented as a union of at most two intervals, can be computed in $O(1)$ time in the PILLAR model.

Proof. First we explain how to compute $\operatorname{PairMatch}(T, P, i, 0)$. Let $p(i) = \text{LCP}(T[i..], P)$ and $s(i) = \text{LCP}_R(T[..i), P)$. If $p(i) + s(i) \geq m$, an interval $[i - s(i) .. i + p(i) - m]$ of starting positions of circular occurrences of $P$ in $T$ is reported; otherwise the answer is an empty set.

In general $\operatorname{PairMatch}(T, P, i, j)$ can be computed using (at most) two queries of the type $\operatorname{PairMatch}(T, P, i', 0)$, for $i' = i - j$ and $i' = i - j + m$. A respective query is asked only if $i' \in [0..n - m)$. The resulting intervals need to be intersected with $(i - m..i]$ to ensure that the circular occurrence contains position $i$.

6.3 Internal CPM

The circular pattern matching problem is formally defined as follows.

<table>
<thead>
<tr>
<th>Circular Pattern Matching (CPM)</th>
</tr>
</thead>
<tbody>
<tr>
<td><strong>Input:</strong> A text $T$ of length $n$ and a pattern $P$ of length $m$.</td>
</tr>
<tr>
<td><strong>Output:</strong> All positions in $T$ where circular occurrences of $P$ start.</td>
</tr>
</tbody>
</table>

We will show an efficient solution in the PILLAR model of CPM in the case when the lengths of the pattern and of the texts are similar. The algorithm below applies the results of [7, 8]. These results considered the approximate CPM problem with $k \geq 1$ mismatches or edits. In the proof of the following theorem, we show that they can be adapted to the case of the exact CPM problem, obtaining an even simpler algorithm. The main idea of the algorithm is illustrated in Figure 7.

▶ Theorem 29. If $n \leq 2m$, the answer to the CPM problem, represented as a union of $O(1)$ interval chains, can be computed in $O(1)$ time in the PILLAR model.

Proof. Let $P = P_1P_2$, where $|P_1| = \lfloor m/2 \rfloor$. Each circular occurrence of $P$ in $T$ implies a standard occurrence of at least one of $P_1$ and $P_2$ in $T$. Henceforth, we assume that it implies an occurrence of $P_1$; the remaining case can be treated symmetrically.

Let $A = \operatorname{Occ}(P_1, T)$. As $|T| \leq 4|P_1| + 3$, a representation of $A$ consisting of $O(1)$ arithmetic sequences can be computed using $O(1)$ IPM queries by the so-called standard trick. We consider each of the arithmetic sequences $B$ separately.
Nonperiodic case. If an arithmetic sequence $B$ contains at most two occurrences, then we ask a query \textsc{PairMatch}(T, P, i, 0) for each $i \in B$. The resulting intervals contain positions of all circular occurrences of $P$ in $T$ that imply an occurrence of $P_1$ in $T$ at a position $i \in B$, and possibly some other circular occurrences of $P$ in $T$ (that imply an occurrence of $P_2$).

Periodic case. Assume now that an arithmetic sequence $B$ contains at least three elements. As already mentioned, its difference is $q := \per(P_1)$.

Let $i$ be any element of $B$. We compute the largest index $i_L < i$ and the smallest index $i_R > i$ such that

$$T[i_L] \neq T[i_L + q] \quad \text{or} \quad i_L = -1, \quad T[i_R] \neq T[i_R - q] \quad \text{or} \quad i_R = |T|.$$

Let $Q = P_2P_1P_2$ and $j = |P_2|$. Similarly (see Figure 5), we compute the largest index $j_L < j$ and the smallest index $j_R > j$ such that

$$Q[j_L] \neq Q[j_L + q] \quad \text{or} \quad j_L = -1, \quad Q[j_R] \neq Q[j_R - q] \quad \text{or} \quad j_R = |Q|.$$

The indices $i_L, i_R, j_L, j_R$, which can be called \textit{misperiods}, can be computed using a constant number of \textsc{LCP} and \textsc{LCP}_R queries on $T$ and $P$.

![Figure 5 Misperiods $i_L, i_R, j_L$; in this case, there is no misperiod $j_R$.](image)

We consider two cases:

Case (1). The cyclic occurrence is an occurrence of a rotation of $P$ that is a length-$m$ substring of $Q(j_L \ldots j_R)$; the occurrence is contained within a substring $T(i_L \ldots i_R)$ in the text. Both strings in scope are periodic with period $q$; it only matters if the periods are synchronized. Let

$$X = (j_L \ldots j_R - m) \quad \text{and} \quad Z = (i_L \ldots i_R - m).$$

The set $X$ consists of the positions in $Q$ where a rotation of $P$ contained in $Q(j_L \ldots j_R)$ starts. The set $Z' := \{z \in Z : \exists x \in X \ z \equiv x \ (\text{mod} \ q)\}$ consists of the starting positions of circular occurrences of $P$ contained in $T(i_L \ldots i_R)$. By the following claim, the set $Z'$ can be computed in $O(1)$ time.

\begin{itemize}
  \item Claim 30 ([7, Lemma 7]). Let $X$ and $Z$ be intervals and $q$ be a positive integer. The set $Z' := \{z \in Z : \exists x \in X \ z \equiv x \ (\text{mod} \ q)\}$, represented as a disjoint union of at most three interval chains, can be computed in $O(1)$ time.
\end{itemize}

Case (2). In this case, two misperiods, one in $T$ and one in $P$, need to be synchronized. It suffices to take the union of results of a \textsc{PairMatch}(T, P, i_L, |P_1| + j_L) query if neither of $i_L, j_L$ equals $-1$ and of a \textsc{PairMatch}(T, P, i_R, j_R - |P_2|) query if $i_R \neq |T|$ and $j_R \neq |P|$.

Overall, the result is a union of $O(1)$ intervals and interval chains and can be computed in $O(1)$ time in the \textsc{PILLAR} model using Lemma 28 and Claim 30.
The interval \(X\) (shaded box) represents the starting positions of the rotations of \(P = (aabaa)^4aa\) contained in \(Q(j_L \ldots j_R) = Q(8 \ldots 33)\). Five copies of \(X\) (two of them partial) constitute the output set \(Z'\) (the shaded boxes in Figure 7).

We introduce the following generalization of IPM queries.

**INTERNAL CIRCULAR PATTERN MATCHING QUERIES (INTERNALCPM)**

**Input:** A string \(S\) of length \(n\).

**Queries:** Given two substrings \(P\) and \(T\) of \(S\) such that \(|T| \leq 2|P|\), report all the starting positions of all circular occurrences of \(P\) in \(T\).

Combining Theorems 27 and 29, we obtain the following result, which generalizes Theorem 6.

**Theorem 31.** The answer to an INTERNALCPM query, represented as a union of \(O(1)\) interval chains, can be computed in \(O(1)\) time after \(O(n)\)-time preprocessing.

### 7 Final Remarks

We took a recursive approach proposed in the computation of seeds and adjusted it to the case of cyclic covers. Despite the similarity, several major changes were necessary due to circularity. We hope that such a recursive approach can be used in other problems on strings.

We also demonstrated the importance of a new tool in computations on cyclic strings: internal circular pattern matching queries. Hopefully, they could be used for other problems related to cyclic substrings.

### References


Faster Prefix-Sorting Algorithms for Deterministic Finite Automata

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Abstract
Sorting is a fundamental algorithmic pre-processing technique which often allows to represent data more compactly and, at the same time, speeds up search queries on it. In this paper, we focus on the well-studied problem of sorting and indexing string sets. Since the introduction of suffix trees in 1973, dozens of suffix sorting algorithms have been described in the literature. In 2017, these techniques were extended to sets of strings described by means of finite automata: the theory of Wheeler graphs [Gagie et al., TCS’17] introduced automata whose states can be totally-sorted according to the co-lexicographic (co-lex in the following) order of the prefixes of words accepted by the automaton. More recently, in [Cotumaccio, Prezza, SODA’21] it was shown how to extend these ideas to arbitrary automata by means of partial co-lex orders. This work showed that a co-lex order of minimum width (thus optimizing search query times) on deterministic finite automata (DFAs) can be computed in $O(m^2 + n^{5/2})$ time, $m$ being the number of transitions and $n$ the number of states of the input DFA.

In this paper, we exhibit new combinatorial properties of the minimum-width co-lex order of DFAs and exploit them to design faster prefix sorting algorithms. In particular, we describe two algorithms sorting arbitrary DFAs in $O(mn)$ and $O(n^2 \log n)$ time, respectively, and an algorithm sorting acyclic DFAs in $O(m \log n)$ time. Within these running times, all algorithms compute also a smallest chain partition of the partial order (required to index the DFA). We present an experiment result to show that an optimized implementation of the $O(n^2 \log n)$-time algorithm exhibits a nearly-linear behaviour on large deterministic pan-genomic graphs and is thus also of practical interest.

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1 Introduction

In this paper, we study the problem of indexing string sets for pattern matching queries: pre-process a set $L \subseteq \Sigma^*$ of strings from a finite alphabet $\Sigma$ so that later we can efficiently answer queries of the form “is a given query pattern $P \in \Sigma^*$ substring of some string in $L$?”

Clearly, an algorithmic solution to this problem requires the set $L$ to be representable in finite space (even though $L$ itself could contain an infinite number of strings); in this paper, we focus on string sets described by finite state automata, that is, on regular languages. Our results build on a successful line of previous research based on the following idea: after sorting all prefixes $\text{Pref}(L)$ of the strings in $L$ in colexicographic (co-lex for brevity) order$^1$, pattern matching queries translate to finding the strings in $\text{Pref}(L)$ that are suffixed by pattern string $P$. Being $\text{Pref}(L)$ co-lex sorted, those strings form a range in co-lex order; notice that, if the sorted $\text{Pref}(L)$ is explicitly stored, such a range can be easily found by binary search. Recall, however, that (due to limited available working space) we work with a particular representation of $L$: a finite state automaton $A$. This requires re-formulating the pattern matching problem on $A$. It is easy to see that pattern matching queries on $L$ translate to finding paths of $A$ whose labels, when concatenated, form $P$. When using $A$ to index $L$, the main question becomes therefore “how does the total co-lex order on $\text{Pref}(L)$ map onto the states of $A$?”. In particular cases, such a mapping yields a total order among $A$’s states. This happens, for example, when $A$ is a path (i.e. a string; corresponding data structures include the suffix tree [22], the suffix array [18, 13], and the FM-index [10]), a finite set of disjoint paths (eBWT [19]), or a labeled arborescence (XBWT [9]). A total order on the states of $A$ is obtained even in particular cases where $A$ may accept an infinite language: this is the case, for example, of de Bruijn graphs (BOSS [2]) and Wheeler graphs [11] (the latter generalize all the above classes of totally-sortable labeled graphs).

More recently, in [6, 5] it was shown that in the general case (arbitrary NFAs) the total co-lex order on $\text{Pref}(L)$ maps very naturally onto a family of partial co-lex orders among the states of $A$. Such a family contains only one order for any given DFA, while NFAs may admit multiple admissible co-lex orders. Letting $p$ be the width of a smallest-width partial order $<_A$, it was shown that pattern matching queries on $L$ can be solved in time $\tilde{O}(p^2)$ per query character$^2$. Note that this generalizes the total order case $p = 1$, where indeed queries take $\tilde{O}(1)$ time using the aforementioned solutions (e.g. indexes on strings and labeled trees). Building the index of [6, 5] requires the computation of a smallest chain partition for the co-lex order $<_A$, i.e. a minimum-size partition $C_1, \ldots, C_p$ of $A$’s states such that $(C_i, <_A)$ is a total order for each $i = 1, \ldots, p$ (note that the index does not require the order $<_A$ itself, just a chain partition). Letting $n$ and $m$ be the number of states and transitions of $A$, respectively, [6] showed how to build such a chain partition in $O(n^{5/2} + m^2)$ time in the case where $A$ is a deterministic finite automaton (DFA). The work [5] presented a solution running in $\tilde{O}(m^2)$ time w.h.p. In the general nondeterministic (NFA) case, the problem is known to be NP-complete$^3$, even though polynomial algorithms do exist for co-lex pre-orders [4] (which still allow indexing and whose width is never larger than that of co-lex orders).

---

1 Historically, the lexicographic order of suffixes was used first; however, with finite state automata the symmetric co-lex order of $\text{Pref}(L)$ turns out to be more natural.

2 The notation $\tilde{O}$ hides factors polylogarithmic in the size of $A$.

3 Hardness follows from hardness of the $p = 1$ case [12], while membership in NP follows from the fact that the properties defining a co-lex order can be checked in polynomial time, given a candidate order.
1.1 Our results

In this work, we focus on the problem of computing the smallest-width partial co-lex order \( <_A \) when the input is a DFA. On DFAs, \( <_A \) has a very intuitive definition: letting \( u, v \) be states of \( A \), we have \( u <_A v \) if and only if \( \alpha < \beta \) for every \( \alpha \in I_u \) and \( \beta \in I_v \), where \( < \) denotes the co-lex order among strings and \( I_u \) denotes the set of strings (in fact, a regular language) labeling all paths from the source of \( A \) to \( u \). We first observe that \( <_A \) is completely specified by pairs \((\inf I_u, \sup I_u)\) over the co-lex sorted \( \text{Pref}(L) \): in fact, we prove that \( u <_A v \) holds if and only if \( \sup I_u \leq \inf I_v \). This allows finding a smallest chain decomposition of \( <_A \) in \( O(n) \) time through a solution of the interval partitioning problem, given that the co-lex ranks of strings \( \inf I_u \) and \( \sup I_u \) are known for each state \( u \). Observing that these strings can be easily encoded with two pruned versions of the DFA \( A \), this leaves the problem of computing and sorting them — ideally, in \( O(m) \) time. We give three different solutions for this problem, which could be of independent interest. The first two solutions work on arbitrary DFAs and run in time \( O(mn) \) and \( O(n^2 \log n) \), respectively. The latter of these two solutions is based on suffix doubling, the technique at the core of the first suffix array construction algorithm [18], and is close to optimal on dense graphs. We show that an optimized implementation of this algorithm exhibits a sub-quadratic behaviour on large deterministic pan-genomic graphs (in fact, we experimentally observe a linearithmic running time). The third solution works on acyclic DFAs, runs in \( O(m \log n) \) time, and generalizes a well-known algorithm for building the Burrows-Wheeler transform in an online fashion; in our case, we process the automaton’s states in any topological order and, for each processed state \( u \), compute \( \inf I_u \) and \( \sup I_u \) using the results computed on the already-processed states.

2 Preliminaries

Notation \([i,j]\), where \( i, j \in \mathbb{N} \), denotes the integer set \( \{i, i+1, \ldots, j\} \) (if \( i > j \), then \([i,j] = \emptyset\)).

Let \( \Sigma \) be a finite alphabet. A finite string \( \alpha \in \Sigma^* \) (or string of finite length) is a finite concatenation of characters from \( \Sigma \). The notation \(|\alpha|\) indicates the length of the string \( \alpha \). The symbol \( \epsilon \) denotes the empty string. The notation \( \alpha[i] \) denotes the \( i \)-th character from the beginning of \( \alpha \); indices start from 1, so \( \alpha[1] \) is the first character of \( \alpha \). Letting \( \alpha, \beta \in \Sigma^* \), \( \alpha \cdot \beta \) (or simply \( \alpha \beta \)) denotes the concatenation of strings. The notation \( \alpha[i..j] \) denotes \( \alpha[i] \cdot \alpha[i+1] \cdot \ldots \cdot \alpha[j] \); if \( i > j \), then \( \alpha[i..j] \) is the empty string \( \epsilon \). The notation \( \alpha \subseteq \beta \), where \( \alpha, \beta \in \Sigma^* \), indicates that \( \alpha \) is a prefix of \( \beta \), i.e. \( \alpha = \beta[1..i] \) for some \( i \leq |\beta| \). An \( \omega \)-string \( \beta \in \Sigma^\omega \) (or infinite string / string of infinite length) is an infinite numerable concatenation of characters from \( \Sigma \). In this paper, we work with left-infinite \( \omega \)-strings, meaning that \( \beta \in \Sigma^\omega \) is constructed from the empty string \( \epsilon \) by prepending an infinite number of characters to it. In particular, the operation of appending a character \( a \in \Sigma \) at the end of a \( \omega \)-string \( \alpha \in \Sigma^\omega \) is well-defined and yields the \( \omega \)-string \( \alpha a \). The notation \( \alpha^\omega \), where \( \alpha \in \Sigma^* \), denotes the concatenation of an infinite (numerable) number of copies of string \( \alpha \).

Definition 1. A Deterministic Finite-State Automaton (DFA) is a quintuple \( A = (Q, \Sigma, \delta, s, F) \) where \( Q \) is the finite set of states, \( \Sigma \) is a finite alphabet, \( \delta : Q \times \Sigma \rightarrow Q \) is the transition function, \( s \in Q \) is the initial state, and \( F \subseteq Q \) is the set of final states.

As is customary, we extend the transition function to words \( \alpha \in \Sigma^* \) as follows: for \( a \in \Sigma \), \( \alpha \in \Sigma^* \), and \( q \in Q \): \( \delta(q, a \cdot \alpha) = \delta(\delta(q, a), \alpha) \) and \( \delta(q, \epsilon) = q \). By \( \delta^{-1}(u) \), we denote the set of states from which there exists a transition to \( u \) i.e. \( \delta^{-1}(u) = \{ v \in Q : (\exists a \in \Sigma)(\delta(v, a) = u) \} \).

In the rest of the paper, \( n = |Q| \) denotes the number of states and \( m = |\delta| = |\{(u, v, a) \in Q \times Q \times \Sigma : \delta(u, a) = v\}| \) the number of transitions of the DFA under consideration.
Following [1], we use the following notation for the set of words reaching a given state:

Definition 2. Let $A = (Q, \Sigma, \delta, s, F)$ be a DFA. If $q \in Q$, let $I_q$ be the set of words reaching $q$ from the initial state:

$$I_q = \{ \alpha \in \Sigma^* : q = \delta(s, \alpha) \};$$

$I_q$ is also called the regular language recognized by $q$.

The language $L(A)$ recognized by $A$ is defined as $L(A) = \cup_{q \in F} I_q$.

The co-lexicographic (or co-lex) order of two strings $\alpha, \beta \in \Sigma^+ \cup \Sigma^*$ is defined as follows.

(i) $\epsilon < \alpha$ for every $\alpha \in \Sigma^+ \cup \Sigma^*$, and (ii) if $\alpha = \alpha'a$ and $\beta = \beta'b$ (with $a, b \in \Sigma$ and $\alpha', \beta' \in \Sigma^* \cup \Sigma^*$), $\alpha < \beta$ holds if and only if $(a < b) \lor (a = b \land \alpha' < \beta')$. In this paper, the symbols $<$ and $\leq$ will be used to denote the total order between the alphabet’s characters, the co-lexicographic order between strings/$\omega$-strings, and the co-lex partial order among the states of an automaton (Definition 3). The meaning of symbols $<$ and $\leq$ will always be clear from the context. In all cases, the symbol $\leq$ has the following meaning: $x \leq y$ if and only if $x < y$ or $x = y$ (i.e., $x < y$ or $x = y$ and $y$ are the same state, the same character, or the same string, depending on the context).

Let $A = (Q, \Sigma, \delta, s, F)$ be a DFA. We assume that $s$ has no incoming edges; any automaton can always be transformed into an equivalent automaton with this property. We also assume that every state is reachable from the source: for every $v \in Q$, there exists $\alpha \in \Sigma^*$ such that $\delta(s, \alpha) = v$. Moreover, we assume input consistency: for every $u, v, v' \in Q$ and $c, c' \in \Sigma$, if $\delta(v, c) = \delta(v', c') = u$, then $c = c'$. We denote with $\lambda(v)$ such a uniquely-defined character and take $\lambda(s) = \#$, for the source $s$, where $\# \notin \Sigma$ is such that $\# < c$ for every $c \in \Sigma$.

Note that input consistency is equivalent to working with state-labeled automata. Also this assumption is not too restrictive, since any automaton can be converted into an equivalent input-consistent automaton by just multiplying its size by a factor of $|\Sigma|$.

The following concepts can be defined more in general for NFAs (see [6]), but for the purposes of this article it will be sufficient to introduce them just on DFAs:

Definition 3. Let $A = (Q, \Sigma, \delta, s, F)$ be a DFA. A co-lex order on $A$ is a partial order $\leq$ on $Q$ that satisfies the following two axioms:

1. (Axiom 1) For every $u, v \in Q$, if $u < v$, then $\lambda(u) \leq \lambda(v)$;
2. (Axiom 2) For every $a \in \Sigma$ and $u, v, u', v' \in Q$, if $u = \delta(u', a)$ and $v = \delta(v', a)$ and $u < v$, then $u' \leq v'$.

The width of a partial order is the size of its largest antichain or, equivalently by Dilworth’s theorem [7], the size of a smallest chain partition of the order.

Definition 4. The co-lex width of a DFA $A$ is the minimum width of a co-lex order on $A$:

$$\text{width}(A) = \min \{ \text{width}(\leq) : \leq \text{ is a co-lex order on } A \}$$

On DFAs, the following co-lex order is of particular interest:

Definition 5. Let $A$ be a DFA. The relation $<_A$ over $Q$ is defined by:

$$u <_A v \text{ if and only if } (\forall \alpha \in I_u)(\forall \beta \in I_v) (\alpha < \beta).$$

In fact, by [5, Lem. 1] the following holds:

Lemma 6. If $A$ is a DFA, then $<_A$ is a co-lex order on $A$ and $\text{width}(<_A) = \text{width}(A)$. The order $<_A$ is called the maximum co-lex order on $A$. 

Computing the smallest-width co-lex order is of interest because, as shown in [5, 6], there exists a linear-space index over any DFA $\mathcal{A}$ answering subpath queries (find all the states of $\mathcal{A}$ reached by a path labeled with a given query string $P$) in time proportional to $\text{width}(\mathcal{A})^2$ time per query character. In fact, the index is even compressed and uses $\log(\text{width}(\mathcal{A})) + \log(\Sigma) + O(1)$ bits per transition of $\mathcal{A}$. Building such an index requires computing a smallest-size chain partition of $<_{\mathcal{A}}$. State-of-the-art algorithms for this problem run in time $O(m^2 + n^{3/2})$ [6] and $\tilde{O}(m^2)$ w.h.p. [5]. The goal of our paper is to improve these bounds by exploiting a new characterization for $<_{\mathcal{A}}$, introduced in the next section.

3 A new characterization of the maximum co-lex order of a DFA

In this section, we give a new interval-based characterization of the maximum co-lex order $<_{\mathcal{A}}$ of a DFA. We show that this yields an $O(n)$-space representation for $<_{\mathcal{A}}$ (observe that, in general, a partial order requires $O(n^2)$ space to be represented) and that, given this representation, one can compute a smallest chain partition of $<_{\mathcal{A}}$ in linear $O(n)$ time.

3.1 Infimum and supremum strings

Let $u$ be a state of a DFA $\mathcal{A} = (Q, \Sigma, \delta, s, F)$. For the set $I_u$ of strings recognized by $u \in Q$, consider a (possibly infinite) string $\beta$ such that $\beta$ is a lower bound of $I_u$; i.e., $\beta \leq \alpha$ for every $\alpha \in I_u$. Consider the co-lex largest string $\gamma$ among such lower bounds of $I_u$. We call such a string the infimum string of $u$, and denote it by $\inf I_u$. Similarly, we define the supremum string $\sup I_u$ of $u$ as the least upper bound of $I_u$; see Figure 1 for an example.

**Definition 7** (Infimum and supremum strings). Let $u \in Q$ be a state of a DFA $\mathcal{A} = (Q, \Sigma, \delta, s, F)$. The infimum string $\inf I_u$ and the supremum string $\sup I_u$ are defined as:

\[
\inf I_u = \gamma \in \Sigma^* \cup \Sigma^\omega \text{ s.t. } (\forall \beta \in \Sigma^* \cup \Sigma^\omega \text{ s.t. } (\forall \alpha \in I_u \beta \leq \alpha) \beta \leq \gamma)
\]

\[
\sup I_u = \gamma \in \Sigma^* \cup \Sigma^\omega \text{ s.t. } (\forall \beta \in \Sigma^* \cup \Sigma^\omega \text{ s.t. } (\forall \alpha \in I_u \alpha \leq \beta) \gamma \leq \beta)
\]

![Figure 1 Example DFA with its infimum/supremum strings.](image)

As a warm up, we make several observations on $I_u$, infimum, and supremum strings.

**Observation 8.** Let $\mathcal{A} = (Q, \Sigma, \delta, s, F)$ be a DFA. For any $u \in Q$, the following hold:

1. For every $\alpha \in I_u$, $\alpha$ is finite.
2. For any $v \neq u \in Q$, $I_u \cap I_v$ is the empty set.
3. For any finite suffix $\alpha \in \Sigma^*$ of $\inf I_u$ (or $\sup I_u$), there exists $\beta \in \Sigma^*$ such that $\beta \alpha \in I_u$.
4. $\inf I_u \in I_u$ if and only if $\inf I_u$ is a finite string; similar for $\sup I_u$.
5. $I_u$ is a singleton if and only if $\inf I_u = \sup I_u$. In such a case, $I_u = \{\inf I_u(= \sup I_u)\}$ and $\inf I_u = \sup I_u \in I_u$ is a finite string.
6. For $v \neq u \in Q$, if $\inf I_u = \inf I_v$ or $\inf I_u = \sup I_v$ then $\inf I_v$ has infinite length; similar for $\sup I_v$.

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Let $\mathcal{A} = (Q, \Sigma, \delta, s, F)$ be a DFA. For any $u \in Q$, the following hold:

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4. $\inf I_u \in I_u$ if and only if $\inf I_u$ is a finite string; similar for $\sup I_u$.
5. $I_u$ is a singleton if and only if $\inf I_u = \sup I_u$. In such a case, $I_u = \{\inf I_u(= \sup I_u)\}$ and $\inf I_u = \sup I_u \in I_u$ is a finite string.
6. For $v \neq u \in Q$, if $\inf I_u = \inf I_v$ or $\inf I_u = \sup I_v$ then $\inf I_v$ has infinite length; similar for $\sup I_v$. 
Proof.
1. By definition of $I_u$.
2. By definition of DFA, for any string $\alpha$, there exists only one state $u$ such that $\delta(s, \alpha) = u$. This will prove our main claim, since for any string $\beta \in I_u$, $\delta(s, \cdot \alpha) = \delta(s, \beta, \alpha)) = \delta(v, \alpha) = u$ by definition, so $\beta \cdot \alpha \in I_u$.
3. Let $\alpha$ be any finite suffix of $\inf I_u$; the sup $I_u$ case is analogous. We claim that there must exist $v \in Q$ such that $\delta(v, \alpha) = u$. This will prove our main claim, since for any string $\beta \in I_u$, $\delta(s, \beta \cdot \alpha) = \delta(s, \beta, \alpha)) = \delta(v, \alpha) = u$ by definition, so $\beta \cdot \alpha \in I_u$.

To prove the claim, assume by contradiction that there is no such $v \in Q$. Let $\alpha' \in \Sigma^*$ be the longest suffix of $\alpha$ such that there exists $v' \in Q$ such that $\delta(v', \alpha') = u$. Let $\alpha'' \in \Sigma^* \cup \Sigma^\omega$ and $a \in \Sigma$ be a string and a symbol such that $\inf I_u = \alpha'' \cdot a \cdot \alpha'$. Let $b \in \Sigma$ be the smallest alphabet symbol that is greater than $a$. Note that such $b$ must exist; if not, every $v'$ must have an incoming transition labeled by a symbol $a' \in \Sigma$ which is co-lex smaller than $a$; since $\alpha' \neq a$ (otherwise $\alpha'$ would be longer), there exists a string in $I_u$ suffixed by $a' \alpha'$ which is co-lex smaller than $\inf I_u$, which causes a contradiction. Then, for every such $v'$ and every $v'' \in Q$ and $c \in \Sigma$ such that $\delta(v'', c) = v'$, we have $a < b \leq c$. Note that $\inf I_u < b \cdot \alpha' \leq \gamma$ for every $\gamma \in I_u$, so $\inf I_u$ is not the greatest lower bound of $I_u$. This is a contradiction with the definition of $\inf I_u$.

4. $(\Rightarrow)$ By definition of $I_u$.

$(\Leftarrow)$ Let $\inf I_u$ be finite. Let us assume, by contradiction, that $\inf I_u \notin I_u$. Let $\alpha = a \cdot \inf I_u$ where $a \in \Sigma$ is the smallest symbol of the alphabet. We claim that $(\inf I_u <) \alpha \leq \beta$ for every $\beta \in I_u$, which contradicts the definition of $\inf I_u$. Consider $\beta \in I_u$. Let $k$ be the length of the longest common suffix of $\beta$ and $\inf I_u$. If $k < |\inf I_u|$, then obviously $\alpha < \beta$ because prepending a symbol to $\inf I_u$ does not affect the relative co-lex order of $\inf I_u$ and $\beta$. If $k = |\inf I_u|$, then $\inf I_u$ is a suffix of $\beta$ and $|\inf I_u| + 1 \leq |\beta|$ because $\inf I_u \notin I_u$.

Therefore after prepending the smallest symbol $a$ to $\inf I_u$, we still have $a \cdot \inf I_u \leq \beta$. To prove the other case for $\sup I_u$, let $\sup I_u$ be finite and let us assume for a contradiction that $\sup I_u \notin I_u$. From (3), there exists $\beta \in \Sigma^*$ such that $\beta \alpha \in I_u$ where $\alpha = \sup I_u$; note that a string is a suffix of itself. Because $\delta(s, \alpha) \neq u$, it holds $\beta \neq \epsilon$. However, then we have $\sup I_u = \alpha < \beta \alpha \in I_u$, which contradicts with the definition of $\sup I_u$ being an upper bound of $I_u$.

5. $(\Rightarrow)$ If $I_u = \{\alpha\}$, then clearly $\inf I_u = \alpha$ and $\sup I_u = \alpha$, so $\inf I_u = \sup I_u$.

$(\Leftarrow)$ If $\inf I_u = \sup I_u$, then they are the same finite string. To see this, assume by contradiction that $\inf I_u = \sup I_u$ have infinite length. Then, for every $\alpha \in I_u$, $\inf I_u < \alpha < \sup I_u$. Since $\inf I_u = \sup I_u$, no such $\alpha$ can exist thus $I_u = \emptyset$. This is a contradiction, because it must hold $I_u \neq \emptyset$ by the assumption that there always exists $\alpha \in \Sigma^*$ such that $\delta(s, \alpha) = u$ (and, for the source $s$, $I_u = \{\epsilon\}$). Since $\inf I_u = \sup I_u$ is a finite string, $\inf I_u \in I_u$ by (4). In addition, $\inf I_u$ is the unique string in $I_u$ because for every $\alpha \in I_u$, $\inf I_u \leq \alpha \leq \sup I_u$ and $\inf I_u = \sup I_u$, therefore $\inf I_u = \alpha = \sup I_u$.

6. Immediate from Observations (2) and (4).

To conclude the section, we prove a lemma showing that infimum and supremum strings can always be expressed as a (possibly, infinite) concatenation of a constant number of distinct strings whose length does not exceed the number of states. This lemma will be useful later to bound the sorting depth of our algorithms computing $\prec_A$.

Lemma 9. For a DFA $A = (Q, \Sigma, \delta, s, F)$ and a state $u \in Q$, let $\gamma \in \{\inf I_u, \sup I_u\}$ be either the infimum or the supremum string of $I_u$. Then,
1. If $\gamma$ is finite, then $|\gamma| < |Q|$.
2. If $\gamma$ has infinite length, then $\gamma = \beta^{*} \alpha$ for some $\alpha, \beta \in \Sigma^*$ such that $|\alpha| + |\beta| < |Q|$.
Proof. Suppose $|\gamma| \geq |Q|$. Let $\gamma = \gamma''$, where $\gamma''$ is the length-$|Q|$ suffix of $\gamma$. Consider a sequence of states $v_1, v_2, \ldots, v_{|Q|+1}$ such that $v_{|Q|+1} = u$ and $\delta(v_k, \gamma''[k]) = v_{k+1}$ for $1 \leq k \leq |Q|$. Note that $v_i \neq s$ for every $2 \leq i \leq |Q| + 1$ because the start state $s$ does not have an incoming transition. Then, there are at most $|Q| - 1$ distinct states among the $|Q|$ states $v_2, \ldots, v_{|Q|+1}$ so by the pigeonhole principle there must be $2 \leq i < j \leq |Q| + 1$ such that $v_i = v_j$. Let $\beta' = \gamma''[i..-1]$, $\beta = \gamma''[i..j-1]$, and $\alpha = \gamma''[j..|Q|]$. Note that $\beta$ is a proper suffix of $\gamma''$ (proper because $i \geq 2$), therefore $|\alpha| + |\beta| = |\beta|$. Note also that (by definition of $\beta'$, $\beta$, and $\alpha$) $\gamma = \beta' \beta \alpha$.

Let us assume $\gamma = \inf I_u$; the other case $\gamma = \sup I_u$ is analogous. Note that $\gamma = \beta' \beta \alpha \leq \beta' \beta^k \alpha$ for every $k \geq 0$. To see this, observe that if $\gamma$ is a finite string, then $\gamma = \beta' \beta \alpha \in I_u$ by Observation 8.4. Since $\delta(s, \beta') = v_1, \delta(v_i, \beta) = v_j = v_i$, and $\delta(v_j = v_i), \alpha = v_{|Q|+1} = u$, we have $\beta' \beta^k \alpha \in I_u$ for every $k \geq 0$. By definition of $\inf I_u$, $\gamma = \inf I_u \leq \beta' \beta^k \alpha$ for every $k \geq 0$.

On the other hand, if $\gamma$ has infinite length, assume for a contradiction that there exists $k' \geq 0$ such that $\beta' \beta^k \alpha < \beta' \beta \alpha = \gamma$. Consider the length-$(l + 1)$ suffix $\beta'' \alpha$ of $\beta' \beta \alpha$ where $l$ is the length of the longest common suffix between $\beta' \beta^k \alpha$ and $\gamma$; clearly, such a $l$ is finite and $l \geq |\alpha|$ since $\alpha$ suffices both strings. Then by Observation 8.3, there exists $\beta'' \beta \alpha \in \Sigma^*$ such that $\beta'' \beta \alpha < \beta' \beta \alpha = \gamma = \inf I_u$, which contradicts the definition of $\inf I_u$.

By plugging $k = 0$ into the inequality $\beta' \beta \alpha \leq \beta' \beta^k \alpha$ above, we obtain $\beta' \beta \alpha \leq \beta' \alpha$. Equivalently (by removing the common suffix $\alpha$) it holds $\beta' \beta \leq \beta'$; but then, we can plug again a common suffix $\beta' \alpha$ for any $k \geq 0$ and obtain that $\beta' \beta^k \alpha \leq \beta' \beta^k \alpha$ for any $k \geq 0$. In particular, this implies that $\beta' \beta^k \alpha = \gamma$ for any $k \geq 1$.

Since in the previous two paragraphs we proved that $\gamma \leq \beta' \beta^k \alpha$ and $\beta' \beta^k \alpha \leq \gamma$ for any $k \geq 1$, we conclude that $\gamma = \beta' \beta \alpha$ for any $k \geq 1$, i.e. $\gamma$ must be an $\omega$-string of the form $\gamma = \beta\omega$. This proves claim (2). Claim (1) also follows since the assumption that $\gamma$ is finite and $|\gamma| \geq |Q|$ leads to $\gamma = \beta\omega$ (a contradiction to the finiteness of $\gamma$), hence its negation (i.e. claim 1) must hold.

### 3.2 $O(n)$-space representation of $<_\mathcal{A}$

Let $\mathcal{K}(\mathcal{A}) = \{\inf I_u : u \in Q\} \cup \{\sup I_u : u \in Q\} \subseteq \Sigma^* \cup \Sigma^\omega$ be the set of all infimum and supremum strings of $\mathcal{A}$. Let $\text{rank}(\alpha)$, for $\alpha \in \mathcal{K}(\mathcal{A})$, denote the position of $\alpha$ in the total order $(\mathcal{K}(\mathcal{A}), <)$ (e.g. $\text{rank}(\alpha) = 1$ for the colex smallest string $\alpha \in \mathcal{K}(\mathcal{A})$, and so on).

Our new representation of $<_\mathcal{A}$ is the set of $n$ integer pairs $\{(\text{rank}(\inf I_u), \text{rank}(\sup I_u)) : u \in Q\} \subseteq [1, 2n] \times [1, 2n]$ (note that $|\mathcal{K}(\mathcal{A})| \leq 2n$). With the next theorem, we show that this set is indeed sufficient to reconstruct $<_\mathcal{A}$.

➤ **Theorem.** Let $\mathcal{A} = (Q, \Sigma, \delta, s, F)$ be a DFA. Then, for any $u, v(\neq u) \in Q$, $u <_\mathcal{A} v$ if and only if $\inf I_u \leq \inf I_v$.

Proof. ($\Rightarrow$) To prove $u <_\mathcal{A} v \Rightarrow \sup I_v \leq \inf I_u$ for all $u, v \in Q$, assume by contradiction that there exist $u, v \in Q$ such that $u <_\mathcal{A} v$ and $\inf I_v < \sup I_u$. We claim that, in this case, there must exist $\alpha \in I_u, \beta \in I_v$ such that $\beta < \alpha$. By Definition 5, this contradicts $u <_\mathcal{A} v$. First, note that there must exist $\alpha \in I_u$ such that $\inf I_v < \alpha$, otherwise it would be $\sup I_v \leq \inf I_v$. We divide the proof by contradiction in the two cases (i) $\inf I_v$ is a finite string and (ii) $\inf I_v$ has infinite length.

(i) If $\inf I_v$ is finite, then $\inf I_v \in I_v$ by Observation 8.4. Choosing $\beta = \inf I_v$, we have $\beta = \inf I_v < \alpha$. This contradicts $u <_\mathcal{A} v$.

(ii) If $\inf I_v$ has infinite length, then by Lemma 9 we can write it as $\inf I_v = \gamma_1 \gamma_2$ for some strings $\gamma_1, \gamma_2 \in \Sigma^*$. Note that, for every $k \geq 0$, there exists a string $\gamma_3 \in \Sigma^*$...
such that $\gamma_3 \gamma_2^k \gamma_1 \in I_v$ (by Observation 8.3 because $\gamma_2^k \gamma_1$ is a suffix of $\inf I_u$). Choose any integer $k'$ such that $|\gamma_2^k \gamma_1| > |\alpha|$ (such an integer exists since $\alpha$ is finite). Since $\inf I_v = \gamma_2^k \gamma_1 < \alpha$, we also have $\gamma_3 \gamma_2^k \gamma_1 (= \beta \in I_v) < \alpha$. Again, this contradicts $u <_A v$.

$(\Leftarrow)$ Let $\sup I_u \leq \inf I_v$, and choose any $\alpha \in I_u$ and $\beta \in I_v$. We need to prove that $\alpha <_A \beta$. By definition of $\sup I_u$ and $\inf I_v$, we have $\alpha \leq \sup I_u \leq \inf I_v \leq \beta$. If $\sup I_u < \inf I_v$, then $\alpha < \beta$. If, on the other hand, $\sup I_u = \inf I_v$ then both $\sup I_u$ and $\inf I_v$ must be finite strings by Observation 8.6. Since $\alpha$ and $\beta$ are both finite, it must be the case that $\alpha \neq \sup I_u$ and $\beta \neq \inf I_v$, therefore $\alpha < \sup I_u = \inf I_v < \beta$. Since this holds for any $\alpha \in I_u$ and $\beta \in I_v$, by definition of $<_A$ it holds $\alpha <_A \beta$. $\blacktriangle$

Equivalently, Theorem 10 shows that $<_A$ can be interpreted as a set of intervals on the co-lex sorted $\text{Pref}(L(A))$. This characterization of $<_A$ will allow us to compute this order faster than the state-of-the-art by (i) co-lex sorting the infimum and supremum strings (Section 4), and (ii) computing a smallest chain partition for $<_A$ in linear time (Section 3.3).

### 3.3 Linear-time chain partitioning algorithm

In general, a partial order over $n$ elements requires $O(n^2)$ space to be represented. Moreover, the fastest general-purpose algorithms for computing the smallest chain partition of a partial order run either in worst-case time $O(n^{5/2})$ (see, for example, [6, Lem. 6.1]) or in $O(n^2)$ time w.h.p. [16]. In this section we show that given the $O(n)$-space representation $S = \{(\text{rank} (\inf I_u), \text{rank} (\sup I_u)) : u \in Q\}$ of $<_A$, from which the order can be represented using intervals, we can compute a smallest chain partition of this order in optimal $O(n)$ time. It is known that the optimal solution of a smallest chain partition of interval orders can be computed with a greedy method (see [14, Sec. 6.8]). Moreover, given the sorted intervals, one can compute it in linear time [3]; for completeness here we give the details.

Based on Theorem 10, we now show a simple linear-time reduction from the smallest chain partition problem (where the input order is represented as described in Section 3.2) to the following problem:

**Definition 11** (Interval partitioning problem, cf. [15, Sec. 4.1]). Let $\{(s_1, f_1), \ldots, (s_n, f_n)\}$ be a set of $n$ activities that must be served (each) by a device. One device can handle at most one activity at the same time. $[s_i, f_i]$ is an interval, where $s_i$ and $f_i$ are the starting and finishing time of activity $i$, respectively. Determine the minimum number of devices to serve all the activities.

Let $S = \{a_1 = (s_1, f_1), \ldots, a_n = (s_n, f_n)\}$ be an instance of the smallest chain partition problem for $<_A$ (that is, a particular instance of $<_A$). Our reduction from this instance to an instance of the interval partitioning problem works as follows:

1. For each pair $a_i = (s_i, f_i)$, with $i \in [1..n]$, let $s_i' = 2s_i + 1$ and $f_i' = 2f_i$.
2. Return the set of intervals $S'' = \{a''_i\}_{i=1}^n$, where $a''_i = [s''_i, f''_i]$, $f''_i = \max(s''_i, f''_i)$

The following Lemma shows that our reduction is correct:

**Lemma 12.** Let $(s_i, f_i), (s_j, f_j)$ be two input pairs, with $s_i \leq s_j$ without loss of generality. Let moreover $[s_i', f_i'], [s_j', f_j']$ be the intervals into which the two pairs get transformed by the above reduction. Then, $f_i \leq s_j$ if and only if $f''_i < s''_j$ (i.e. $[s''_i, f''_i]$ and $[s''_j, f''_j]$ do not overlap).

**Proof.** We divide the proof into two cases: (Case 1) at least one of $s_i = f_i$ or $s_j = f_j$ holds, and (Case 2) both $s_i < f_i$ and $s_j < f_j$ hold.
(Case 1). First, we show that \( f_i \neq s_j \). Assume that \( s_i = f_i \) (the other case \( s_j = f_j \) is analogous). Let \( u \) be the state associated with the pair \((s_i, f_i)\), and \( v \) be the state associated with the pair \((s_j, f_j)\). By Observation 8.5, \( s_i = f_i \) implies that \( \inf I_u = \sup I_u = s_i \). If \( \inf I_v \) is an infinite string, then clearly \( \sup I_v \neq \inf I_v \) (being \( \sup I_u \) a finite string), i.e. \( f_i \neq s_j \). If, on the other hand, \( \inf I_v \) is a finite string, then by Observation 8.4 we have \( \inf I_v \in I_v \); since by Observation 8.2, we have \( I_u \cap I_v = \emptyset \), also in this case we derive that \( \sup I_u \neq \inf I_v \), i.e. \( f_i \neq s_j \).

Knowing \( f_i \neq s_j \), we obtain that \( f_i \leq s_j \) \( \Leftrightarrow \) \( f_i < s_j \) \( \Leftrightarrow \) \( f_i + 1 \leq s_j \). Note that, since \( s_i' = 2s_j + 1 > 2s_j \) (if \( s_j \neq f_j \)) or \( s_i' = 2f_j = 2s_j \) (if \( s_j = f_j \)), we have \( 2s_j \leq s_i'' \). Similarly, \( f_i'' = 2f_i + 1 \). Hence \( 2s_i \leq s_i'' < f_i'' \leq 2f_i + 1 \) (note that \( s_i'' < f_i'' \) always holds for any interval in our reduction). Therefore, we have \( f_i \leq s_j \) \( \Rightarrow \) \( f_i + 1 \leq s_j \) \( \Rightarrow \) \( f_i'' \leq 2f_i + 1 < 2(f_i + 1) \leq 2s_j \leq s_i'' \). For the other direction, note that \( 2f_i \leq f_i'' \) and \( s_i'' \leq 2s_j + 1 \). Then, using these inequalities we obtain: \( f_i'' < s_i'' \Rightarrow 2f_i \leq f_i'' < s_i'' \leq 2s_j + 1 \Rightarrow 2f_i < 2s_j + 1 \Rightarrow f_i \leq s_j \).

(Case 2). In this case, we have \( f_i \leq s_j \) \( \Rightarrow \) \( f_i'' = 2f_i < 2s_j + 1 = s_i'' \Rightarrow f_i'' < s_i'' \). For the other direction, note that \( f_i'' < s_i'' \Rightarrow 2f_i = f_i'' < s_i'' \leq 2s_j + 1 \Rightarrow f_i \leq s_j \).

By Lemma 12, we can now solve smallest chain partition problem for the particular order \( <_A \) by reducing it to an instance of the interval partitioning problem. Moreover, it is easy to see that the reduction works in linear time so the linearity of our strategy relies on the cost of the algorithm we use to solve the latter problem. We can use a greedy method (cf. [3, 8]) to optimally solve the interval partitioning problem (namely, using the smallest possible number of devices). The algorithm processes the intervals in non-decreasing order of starting times, breaking ties arbitrarily. For each interval, we choose any idle device among the available ones. We can keep track of the list of the available devices if the starting and finishing times of the intervals are already sorted. If all devices are busy, we add a new device.

The above-sketched algorithm spends amortized constant time on every activity, plus the time required to sort the input set of intervals. As said earlier, the elements of our input pairs (i.e. before the reduction) are integer values in the range \([1, 2n]\). After the reduction, this range gets expanded to \([2, 4n + 1]\). This allows us to radix-sort the intervals in \( O(n) \) time. As a result, in our scenario we can solve the interval partition problem in \( O(n) \) time and, in particular, find the smallest chain partition of \( <_A \) given its ranked-pair representation in linear time.

4 Co-lex sorting infimum/supremum strings

In this section, we present three algorithms to compute and sort the set containing all infimum and supremum strings of a DFA. The first two algorithms sort the strings in such a way that for every iteration the strings are co-lex sorted with respect to a longer suffix; we present one simple solution that increases the suffix length by 1 at each iteration, and one that doubles the suffix length at each iteration. The third algorithm is a generalization of online BWT construction and is based on the online algorithm for sorting Wheeler DFA presented in [1, Sec. 3.2]. This algorithm works only on acyclic DFAs but has a lower time complexity than the former two solutions.

For ease of explanation, we consider only infimum strings since the supremum string case is analogous. Indeed, one can easily compute and sort both infimum and supremum strings at the same time by creating two copies of the input DFA and then running our algorithms on the union of the two DFAs, extracting the infimum strings on one DFA and the supremum strings on the other DFA while at the same time sorting the union of these two string sets.
4.1 Simple $O(mn)$-time algorithm

Let us establish some notations before describing the algorithm. For a (possibly infinite) string $\alpha$ and an integer $k \geq 0$, we denote by $\text{suf}_k(\alpha)$ the length-$k$ suffix of $\alpha$. When $|\alpha| < k$, we pad $\text{suf}_k(\alpha)$ by prepending $k - |\alpha|$ copies of a special symbol $\# \notin \Sigma$, with $\# < c$ for all $c \in \Sigma$; in this way, we guarantee that $\text{suf}_k(\alpha)$ is always a string of length $k$ and we do not affect the co-lex order of such suffixes (which remains the same before and after the padding).

For state $u \in Q$ and integer $k \geq 0$, we denote by $\text{rank}_k(u) \in \mathbb{N}$ the intermediate rank at iteration $k$ of the co-lex order we are computing; this integer indicates the co-lex rank of $\text{suf}_k(\inf I_u)$ among $\{\text{suf}_k(\inf I_v) : v \in Q\}$. More formally, for $u \in Q$ and $k \geq 0$,

\[
\begin{align*}
\text{rank}_0(u) &= 1 \\
\text{rank}_k(u) &= |\{\text{suf}_k(\inf I_v) : v \in Q \land \text{suf}_k(\inf I_v) \leq \text{suf}_k(\inf I_u)\}| \quad \text{for } k > 0
\end{align*}
\]

Observe that two states are assigned the same rank if their corresponding length-$k$ suffixes are equal. The algorithm works by pruning transitions of the input automaton, i.e. by removing, for every state $u$, transitions coming from a state with a non-minimum rank among the predecessors of $u$. We denote by $\delta_k$ the (pruned) transition function at iteration $k$.

The algorithm works as follows. At iteration $k \geq 0$, we perform the following operations:

1. **Compute $\text{rank}_{k+1}$**. Sort the states $\{u \in Q\}$ by their label $\lambda(u)$ with ties broken by $\text{rank}_k(u)$ for any $v \in \delta_k^{-1}(u)$ (the step below will guarantee that all predecessors $v$ of $u$ have the same $\text{rank}_k(v)$).

2. **Compute $\delta_{k+1}$**. For each $u \in Q$, keep only the transitions from the min-rank predecessors: for $v \in \delta_k^{-1}(u)$, $v \in \delta_{k+1}^{-1}(u)$ iff $\text{rank}_{k+1}(v) = \min\{\text{rank}_{k+1}(u') : u' \in \delta_k^{-1}(u)\}$.

As far as the running time of each iteration is concerned, computing $\text{rank}_{k+1}$ can be done in $O(n)$ time by 2-pass radix sorting (that is, by incoming label and breaking ties by any predecessor’s rank $\text{rank}_k$). Computing $\delta_{k+1}$ takes $O(|\delta_k|) = O(|\delta|) = O(m)$ time. Hence, each iteration takes $O(m)$ time.

Since $\forall k \geq 0$, $\forall u \in Q$, and $\forall v \in \delta_k^{-1}(u)$ we have $\text{suf}_{k+1}(\inf I_u) = \text{suf}_k(\inf I_u) \cdot \lambda(u)$, it is easy to see that the following invariant always holds at the beginning of iteration $k$: the infimum strings are sorted with respect to the co-lex order of their length-$k$ suffixes. This invariant shows that the number of iterations we have to perform is exactly the length of the suffixes that need to be sorted to obtain the correct co-lex order of the infimum strings. We are left to find an upper bound to this length; observe that this is not a trivial problem, since infimum strings may have infinite length.

Consider any two infimum strings $\alpha, \beta \in \{\inf I_u : u \in Q\}$. The upper bound above can be computed by upper-bounding the length of the longest common suffix between $\alpha$ and $\beta$. If any of the two strings is finite, then by Lemma 9 their longest common suffix does not exceed length $n$. If both strings are infinite, then by Lemma 9 we can write them as $\alpha = \alpha_2^* \alpha_1$ and $\beta = \beta_2^* \beta_1$ and we can use the following:

\[\textbf{Lemma 13} \text{ (cf. [19]). For two infinite strings } \alpha = \alpha_2^* \alpha_1 \text{ and } \beta = \beta_2^* \beta_1, \text{ where } \alpha_1, \beta_1 \in \Sigma^* \text{ and } \alpha_2, \beta_2 \in \Sigma^+, \text{ let } \alpha' \text{ and } \beta' \text{ be their suffixes of length } k = |\alpha_2| + |\beta_2| + \max\{|\alpha_1|, |\beta_1|\}. \text{ Then, } \alpha' < \beta' \text{ if and only if } \alpha < \beta.\]

\textbf{Proof.} Without loss of generality, let us assume $|\alpha_1| \leq |\beta_1|$. Moreover, note that without loss of generality we can also assume that $|\alpha_2| + |\alpha_1| > |\beta_1|$; if this does not hold, then re-write $\alpha_1 \leftarrow \alpha_2^* \alpha_1$ for the only integer $k > 0$ such that $|\alpha_1| \leq |\beta_1| < |\alpha_2| + |\alpha_1|$ holds; after the transformation, $\alpha$ can still be written as $\alpha = \alpha_2^* \alpha_1$.\]
If $\alpha_1$ is not a suffix of $\beta_1$, then clearly the longest common suffix between $\alpha$ and $\beta$ is at most $|\alpha_1|$, so the claim holds. Let us assume therefore that $\alpha_1$ is a suffix of $\beta_1$, i.e. $\beta_1 = \beta'_1 \alpha_1$ for some $\beta'_1 \in \Sigma^*$. Since by assumption $|\alpha_2| + |\alpha_1| > |\beta_1|$, note that $|\alpha_2| > |\beta'_1|$. Similarly as above, if $\beta'_1$ does not suffix $\alpha_2$ (i.e. $\beta_1 = \beta'_1 \alpha_1$ does not suffix $\alpha = \alpha_2^* \alpha_1$) then the longest common suffix between $\alpha$ and $\beta$ is at most $|\beta_1|$ and the claim holds, so let us assume that $\beta'_1$ is a suffix of $\alpha_2$, i.e. $\alpha_2 = \alpha_2' \beta'_1$ for some $\alpha_2' \in \Sigma^*$. Since $\alpha_2^* = (\alpha_2')^* = (\beta'_1 \alpha_1)^*$, we conclude that comparing co-lexicographically $\alpha = (\beta'_1 \alpha_1)^* \beta'_1 \alpha_1$ and $\beta = \beta'_1 \beta_1 \alpha_1$ reduces to comparing $(\beta'_1 \alpha_1)^* \beta'_1$ and $\beta_1$. According to [19, Proposition 5], given any $\gamma_1, \gamma_2 \in \Sigma^+$ it is sufficient to compare the length-$k'$ suffixes of $\gamma_1^*$ and $\gamma_2^*$ to determine their co-lex order, where $k' = |\gamma_1| + |\gamma_2| - \gcd(|\gamma_1|, |\gamma_2|)$. Our claim easily follows since $|\beta'_1 \alpha_2'| = |\alpha_2'| = |\alpha_2|$. ▶

**Corollary 14.** The co-lex order of the infimum and supremum strings of a DFA is the same as the co-lex order of their length-$2n$ suffixes.

**Proof.** By Lemma 9, we can represent two infimum/infimum strings $\alpha, \beta$ as $\alpha = \alpha_2^* \alpha_1$ and $\beta = \beta_2 \beta_1$. By the same lemma, each of $|\alpha_1|, |\alpha_2|, |\beta_1|$ and $|\beta_2|$, as well as $|\alpha_1| + |\alpha_2|$ and $|\beta_1| + |\beta_2|$, are bounded by $n$. From Lemma 13, it is sufficient to compare the suffixes of length at most $|\alpha_2| + |\beta_2| + \max\{|\alpha_1|, |\beta_1|\} = \max\{||\alpha_1| + |\alpha_2|\} + |\beta_2|, |\alpha_2| + |\beta_1| + |\beta_2|\})$ of $\alpha$ and $\beta$ in order to discover their co-lex order. Therefore, $2n$ is a sufficient suffix length for sorting all the infimum strings correctly. ▶

Putting everything together, we conclude:

**Lemma 15.** The infimum and supremum strings of an input-consistent DFA $A = (Q, \Sigma, \delta, s, F)$ can be computed and sorted in $O(mn)$ time, where $n = |Q|$ is the number of states and $m = |\delta|$ is the number of transitions.

Equivalently, the above lemma shows that the representation of $<_A$ of Section 3.2 can be computed in $O(mn)$ time. Plugging the linear-time chain partition algorithm of Section 3.3, we obtain:

**Theorem 16.** Given an input-consistent DFA $A = (Q, \Sigma, \delta, s, F)$, we can compute a minimum-size chain partition of $<_A$ in $O(mn)$ time, where $n = |Q|$ is the number of states and $m = |\delta|$ is the number of transitions.

### 4.2 $O(n^2 \log n)$-time suffix doubling algorithm

Instead of increasing the length of the sorted suffixes only by 1 at every iteration, we can double it via a generalization of the prefix doubling algorithm [18], the first suffix array construction algorithm that appeared in the literature. Again, for simplicity we describe the algorithm just for infimum strings; it is easy to modify it so that it computes and sorts the union of all infimum and supremum strings.

Algorithm 1 describes our sorting procedure, which we explain in detail in the rest of the section. At every iteration $k \geq 0$, Algorithm 1 keeps the infimum strings sorted by their length-$2^k$ suffixes. Suppose we already sorted the infimum strings with respect to their length-$2^k$ suffixes. To enable the doubling procedure, we need to show how to compute the length-$2^{k+1}$ suffix of each $\inf I_u$ given as input the length-$2^k$ suffixes of each $\inf I_u$, for all $u \in Q$. Given that the infimum strings are sorted by their length-$2^k$ suffixes, for each $u \in Q$ we can achieve this goal by finding a state $v \in Q$ such that

$$\text{sup}_{2k+1}(\inf I_u) = \text{sup}_{2k}(\inf I_v) \cdot \text{sup}_{2k}(\inf I_u).$$
Faster Prefix-Sorting Algorithms for Deterministic Finite Automata

We call such an extender of \( u \) (at distance \( 2^k \)). More formally, the set \( P_k(u) \) of all extenders of \( u \) at distance \( 2^k \) is defined as:

\[
P_0(u) = \{ v \in \delta^{-1}(u) : (\forall v' \in \delta^{-1}(u))(\lambda(v) \leq \lambda(v')) \}
\]

\[
P_k(u) = \{ v \in Q : \delta(v, \text{sf}_{2k}(\inf I_u)) = u \land \text{sf}_{2k}(\inf I_v) \subseteq \text{sf}_{2k+1}(\inf I_u) \} \quad \text{for } k > 0
\]

For \( u \in Q \), let \( \text{rank}_{2k}(u) \) be the co-lex rank of \( \text{sf}_{2k}(\inf I_u) \), as defined in the previous section. Observe that, by definition, for every \( u \in Q \) and \( v_1, v_2 \in P_k(u) \), \( \text{rank}_{2k}(v_1) = \text{rank}_{2k}(v_2) \).

We implement a suffix doubling step as follows. Assume \( P_k(u) \) and \( \text{rank}_{2k}(u) \) have been computed for all \( u \in Q \). We associate with \( u \) the pair \((a_u, b_u)\) where \( a_u = \text{rank}_{2k}(u) \) and \( b_u \) is chosen as follows. If \( P_k(u) \neq \emptyset \), \( b_u = \text{rank}_{2k}(v) \) with any \( v \in P_k(u) \); otherwise, \( b_u = -\infty \) is chosen. Finally, we compute \( \text{rank}_{2k+1}() \) by radix-sorting pairs \((a_u, b_u)\) in \( O(n) \) time.

After computing \( \text{rank}_{2k+1}() \), we need to compute \( P_{k+1}(\cdot) \) for the next doubling step. For a state \( u \in Q \), let \( \hat{P}_{k+1}(u) = \bigcup_{u' \in P_k(u)} P_k(u') \) be the union of the extender sets of \( u \)'s extenders at distance \( 2^k \). Then, we claim that we can compute \( P_{k+1}(u) \) by removing all non-minimum-rank states (i.e. non-minimum \( \text{rank}_{2k+1}() \)) from \( \hat{P}_{k+1}(u) \). The correctness of this procedure follows from the fact that \( P_{k+1}(u) \) can also be defined as the largest subset of \( \hat{P}_{k+1}(u) \) such that, for every \( v \in P_{k+1}(u) \) and \( \hat{v} \in \hat{P}_{k+1}(u) \), \( \text{rank}_{2k+1}(v) \leq \text{rank}_{2k+1}(\hat{v}) \). To see this, first observe that \( P_{k+1}(u) \subseteq \hat{P}_{k+1}(u) \) because (i) \( v \in \hat{P}_{k+1}(u) \)

---

\[^4\] In this case, there are no more characters to be prepended to \( \inf I_u \) (i.e. \( |\inf I_u| < 2^k \)). Since we radix-sort pairs \((a_u, b_u)\), this choice is consistent with the fact that \( \text{sf}_{2k}(\inf I_u) \) is left-padded with copies of symbol \# in order to reach length \( 2^k \), with \# < \( c \) for all \( c \in \Sigma \) (see definition of \( \text{sf}_{2k} \) in the previous section).
Figure 2 The DFA that has a quadratic number of extenders: the \( \sigma = \Theta(n) \) states in the rightmost column (indicated with A) have \( \sigma = \Theta(n) \) extenders each (indicated with B) at distance \( 2^k \), where \( k = 1 \).

If and only if \( \delta(v, suf_{2k+1}(\inf I_u)) = u \), and (ii) \( v \in P_{k+1}(u) \Rightarrow \delta(v, suf_{2k+1}(\inf I_u)) = u \).

Also, by the definition of \( P_{k+1}, suf_{2k+1}(\inf I_u) \) for \( v \in P_{k+1}(u) \) must not be greater than \( suf_{2k+1}(\inf I_u) \) for any \( \tilde{v} \in P_{k+1}(u) \), which is equivalent to \( rank_{2k+1}(v) \leq rank_{2k+1}(\tilde{v}) \); otherwise, \( suf_{2k+1}(\inf I_u) \cdot suf_{2k+1}(\inf I_u) < suf_{2k+1}(\inf I_u) \cdot suf_{2k+1}(\inf I_u) = suf_{2k+2}(\inf I_u) \), which contradicts the definition of \( \inf I_u \).

Since \( rank_{2k+1}(v) \) has already been computed for all \( v \in Q \) and can thus be evaluated in constant time, from the above characterization of \( P_{k+1}(u) \) we obtain that the time required to compute this set is proportional to the time we spend to compute the union \( \hat{P}_{k+1}(u) = \bigcup_{v' \in P_k(u)} P_k(v') \). Observe that, if there were repeated states among the sets \( P_k(u) \), for \( u' \in P_k(u) \), then computing such a union could take time \( O(n^2) \) (for every \( u \in Q \)), leading to a cubic algorithm. Luckily, with the next lemma we show that this is not the case: being the input automaton deterministic, those sets are pairwise disjoint and their union can thus be computed by just concatenating them.

Lemma 17. Let \( u \in Q \) be a state of a DFA, and let \( v_1, v_2(\neq v_1) \in P_k(u) \) be extenders of \( u \) at distance \( 2^k \). Then \( P_k(v_1) \cap P_k(v_2) = \emptyset \).

Proof. Let \( v_1, v_2(\neq v_1) \in P_k(u) \) be extenders of the same state \( u \in Q \) at distance \( 2^k \). Let \( \alpha_1 = suf_{2k}(\inf I_{v_1}) \) and \( \alpha_2 = suf_{2k}(\inf I_{v_2}) \). Assume, for a contradiction, that there exists \( v' \in P_k(v_1) \cap P_k(v_2) \). By definition of \( P_k \), since \( v' \in P_k(v_1) \), it holds \( \delta(v', \alpha_1) = v_1 \). Similarly, it also holds \( \delta(v', \alpha_2) = v_2 \). Since \( v_1, v_2 \in P_k(u) \) are extenders of the same state \( u \in Q \) at distance \( 2^k \), both \( \alpha_1 \) and \( \alpha_2 \) are equal to the length-\( 2^k \) prefix of \( suf_{2k+1}(\inf I_u) \), therefore \( \alpha_1 = \alpha_2 \). Consequently, we have \( v_1 = \delta(v', \alpha_1) = \delta(v', \alpha_2) = v_2 \), i.e., reading a string \( \alpha_1 = \alpha_2 \) from a state \( v' \) we reach two distinct states \( v_1 \neq v_2 \). This is a contradiction with the fact that the automaton is deterministic, so the claim \( P_k(v_1) \cap P_k(v_2) = \emptyset \) must be true.

From Lemma 17, we can compute \( \hat{P}_{k+1}(u) \) in time proportional to its cardinality \( |\hat{P}_{k+1}(u)| \leq n \). Since finding the minimum-rank states can be done in linear \( O(|\hat{P}_{k+1}(u)|) \) time as well, the computation of \( P_{k+1}(u) \) takes time \( O(n) \) for each \( u \in Q \). We conclude that each iteration of the suffix doubling algorithm takes \( O(n^2) \) time.

It is worth noting that, since we keep track of each set \( P_k(u) \), the running time of an iteration is lower-bounded by the total number of extenders therein. In the worst case, however, the total number of extenders at a single iteration could be truly quadratic even on acyclic DFAs: see Figure 2; in this example, there are \( n = 4\sigma + 2 \) states and \( \sigma^2 + 2\sigma + 1 = \Theta(n^2) \) extenders at distance \( 2^k \) for \( k = 1 \).

Putting all together, we have the following result for the suffix doubling algorithm described in this section.
Lemma 18. The infimum and supremum strings of an input-consistent DFA $A = (Q, \Sigma, \delta, s, F)$ can be computed and sorted in $O(n^2 \log n)$ time, where $n = |Q|$ is the number of states.

Equivalently, the above lemma shows that the representation of $<_A$ of Section 3.2 can be computed in $O(n^2 \log n)$ time. Plugging the linear-time chain partition algorithm of Section 3.3, we obtain:

Theorem 19. Given an input-consistent DFA $A = (Q, \Sigma, \delta, s, F)$, we can compute a minimum-size chain partition of $<_A$ in $O(n^2 \log n)$ time, where $n = |Q|$ is the number of states.

The suffix doubling algorithm in practice. Although every iteration of the suffix doubling algorithm needs to keep track of $O(n^2)$ extenders per iteration in the worst case, we conjecture that it is not likely to have a quadratic number of extenders on realistic datasets. To demonstrate this, we conducted a brief experiment using a pan-genomic graph, which is considered to be one of the most important real-world applications of our problem.

We downloaded the Chromosome 22 sequence of the GRCh38 human reference genome and its variation data from 1000 Genome project [17]. This variation dataset contains a set of substitutions, insertions and deletions appearing on the reference human genome sequence collected from 2,548 samples. Using this dataset, we constructed a pan-genomic graph using VG [21], then converted it into a DFA using the classical powerset construction algorithm [20]. We ran an implementation of our suffix doubling algorithm to sort the infimum and supremum strings and measured the number of extenders at each iteration. The largest $P_k(u)$ and $P_k(u)$ (extenders at distance $2^k$ before/after filtering non-minimum-rank states) during the procedure had cardinality 60 and 37, respectively, which might be considered not negligible but quite small when compared to the DFA’s size ($n=51,904,782, m =53,049,316$). In addition, the sum $\sum_{u\in Q} |\hat{P}_k(u)|$ of the number of extender candidates (the union of extenders before filtering non-minimum-rank states) at any fixed distance $2^k$ was at most two times the number of edges, suggesting that in practice our algorithm exhibits a linearithmic complexity on pan-genomic graphs. C++ source code is available at: https://github.com/regindex/DFA-suffix-doubling.

4.3 $O(m \log n)$-time algorithm for acyclic DFAs

If the input DFA is acyclic, then we can sort the infimum strings more efficiently using the algorithm described in [1, Sec. 3.2]. This algorithm processes the states of any acyclic Wheeler DFA $A$ (that is, $\text{width}(A) = 1$) and their incoming edges in any topological order $u_1, \ldots, u_n$ while updating $<_A$ in an online fashion; more precisely, as soon as step 1 $\leq i \leq n$ has finished, the algorithm has computed the total order $<_A$ of the set $\{u_1, \ldots, u_i\}$. The basic idea is to process the states in any topological order while maintaining a dynamic data structure that stores the relative co-lex order of the states according to any representative of $I_u$ (in fact, [1] proves that on Wheeler DFAs, any string in $I_u$ can be chosen as a representative of the whole $I_u$ to sort the automaton’s states). This is possible because, when $u_i \in Q$ is being processed, the structure is able to check if $\text{rank}(v) \leq \text{rank}(u_j)$ for any $v \in \delta^{-1}(u_i)$ and $j < i$, where $\text{rank}(v)$ denotes the position of $v$ in the total order $<_A$ of the already-processed states $u_1, \ldots, u_{i-1}$ (and, by definition of topological order, $\text{rank}(v)$ and $\text{rank}(u_j)$ have already been computed in the previous steps). This information is sufficient to compute $\text{rank}(u_i)$ among the sorted $u_1, \ldots, u_i$.  


In our case (arbitrary acyclic DFAs), we use the above data structure as follows: after topologically sorting \( A \) (in linear time) we process states in this order. When processing state \( u_i \), assume that states \( u_1, \ldots, u_{i-1} \) have already been co-lex sorted according to their strings \( \inf I_{u_1}, \ldots, \inf I_{u_{i-1}} \), using the data structure of [1, Section 3.2]. By scanning the predecessors of \( u_i \), we find the min-rank state \( v^* = \arg \min_{v \in \delta^{-1}(u_i)} \text{rank}(v) \) among them. At this point, we insert state \( u_i \), as well as transition (labeled edge) \( (v^*, u_i, \lambda(u_i)) \), in the data structure. Note that, since only \( (v^*, u_i, \lambda(u_i)) \) is inserted, the data structure of [1, Section 3.2] maintains a spanning tree of \( A \) rooted at the start state \( s \). By construction, it is easy to see that the unique path connecting \( s \) to \( u_i \) in this spanning tree is labeled with string \( \inf I_{u_i} \); the spanning tree encodes the infimum strings. As a result, our sorting problem is equivalent to sorting this spanning tree. It is known that a labeled tree is a special case of Wheeler graphs (see [11]), so the computed co-lex node order of this spanning tree is precisely the co-lex order of all infimum strings.

Since it is immediate to extend the idea to the union of all infimum and supremum strings, taking into account the cost of each update of the data structure [1, Sec. 3.2], we obtain:

**Lemma 20.** The infimum and supremum strings of an input-consistent acyclic DFA \( A = (Q, \Sigma, \delta, s, F) \) can be computed and sorted in \( O(m \log n) \) time where \( n = |Q| \) is the number of states and \( m = |\delta| \) is the number of transitions.

**Proof.** First of all, the data structure [1, Sec. 3.2] supports the following two operations in \( O(\log n) \) time: (i) computing the relative rank of a state among those that are already processed, and (ii) inserting a new edge (a state is inserted into the structure after all its incoming edges have been inserted). As a result, finding \( v^* = \arg \min_{v \in \delta^{-1}(u_i)} \text{rank}(v) \) takes time \( O(|\delta^{-1}(u_i)| \cdot \log n) \). After \( v^* \) has been found, inserting the labeled edge \( (v^*, u_i, \lambda(u_i)) \), as well as state \( u_i \), into the structure takes time \( O(\log n) \). Overall, after all states have been processed the cost of the above operations amounts to \( O(m \log n) \) time. Since a topological order of \( A \) can be computed in \( O(m) \) time, the total running time is \( O(m \log n) \).

Equivalently, the above lemma shows that the representation of \( \prec_A \) of Section 3.2 can be computed in \( O(m \log n) \) time when \( A \) is acyclic. Plugging the linear-time chain partition algorithm of Section 3.3, we obtain:

**Theorem 21.** Given an input-consistent acyclic DFA \( A = (Q, \Sigma, \delta, s, F) \), we can compute a minimum-size chain partition of \( \prec_A \) in \( O(m \log n) \) time, where \( n = |Q| \) is the number of states and \( m = |\delta| \) is the number of transitions.

**References**


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5 In fact, the running time per operation is \( O(\log m') \), where \( m' \) is the number of edges. However, in our case, the number of transitions inserted into the data structure is only \( m' = n - 1 \) since we insert one transition per state (except \( s \)). Therefore, in our case, the time taken per operation is \( O(\log n) \).


Ernesto Lowy-Gallego, Susan Fairley, Xiangqun Zheng-Bradley, Magali Ruffier, Laura Clarke, Paul Fliceck, and The 1000 Genomes Project Consortium. Variant calling on the GRCn38 assembly with the data from phase three of the 1000 Genomes Project. Wellcome Open Research, 4(50), 2020. doi:10.12688/wellcomeopenres.15126.2.


Encoding Hard String Problems with Answer Set Programming

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Abstract
Despite the simple, one-dimensional nature of strings, several computationally hard problems on strings are known. Tackling hard problems beyond sizes of toy instances with straight-forward solutions is infeasible. To solve these problems on datasets of even small sizes, effort has to be put into the conception of algorithms leveraging profound characteristics of the input. Finding these characteristics can be eased by rapidly creating and evaluating prototypes of new concepts in how to tackle hard problems. Such a rapid-prototyping method for hard problems is answer set programming (ASP). In this light, we study the application of ASP on five NP-hard optimization problems in the field of strings. We provide MAX-SAT and ASP encodings, and empirically reason about the merits and flaws when working with ASP solvers.

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1 Introduction

Despite the fact that most string problems found in literature are solvable in polynomial time or even close to linear time or beyond, there are several problems that are known to be NP-hard. Among those, we focus on five problems that are well-perceived regarding the number of publications studying these problems: CLOSEST STRING (csp)\(^1\), CLOSEST SUBSTRING (css), LONGEST COMMON SUBSEQUENCE (lcs), MINIMUM COMMON STRING PARTITION (mcsp), and SHORTEST COMMON SUPERSTRING (scs). These problems have been studied under various viewpoints. With respect to fixed-parameter tractability (FPT), Bulteau et al. [9] gave a comprehensive survey on various NP-hard problems related to strings; this survey comprises the problems studied in this paper. Also, Basavaraju et al. [2] studied the kernelization of a majority of our problems. We address other related work in the individual sections of each problem, but omit references to approximation algorithms due to their amount, and because we put focus on the exact solution of the aforementioned problems formulated as optimization problems.

\(^1\) We stick to the commonly used abbreviation csp in literature despite that cs would fit better with the abbreviations of the other problems.
A major problem in tackling these problems in practice is that naive solutions quickly become impractical with respect to the time complexity. Tailored algorithms\footnote{Meaning that such algorithms usually are based on theoretical results that can be put hardly into practice.} are hard to implement, and thus a burden on the algorithm engineering side. Our contribution is to advertise answer set programming (ASP) as a rapid-prototype programming tool for solving NP-hard string problems on small instances. ASP is a declarative programming language geared towards solving hard problems \cite{40, 12}. ASP has been successfully applied in robotics \cite{3}, or for computing the $n$-queens and the knight’s tour problem \cite{18}. There is also a competition on ASP solvers on various classic problems addressing mainly problems on graphs \cite{28}. See \cite{19, 20} and the references therein for an overview of other use cases.

Although well-devised algorithms can outperform ASP-based approaches, the programming effort for writing in an expressive, declarative programming language such as ASP is considerably small. In this paper, we devise MAX-SAT encodings for the above addressed problems, and subsequently translate these encodings into the ASP language. With respect to tackling hard string problems via MAX-SAT encodings we are aware of the work of Bannai et al. \cite{1} who studied MAX-SAT encodings for repetitiveness measures that are also known to be NP-hard.

## Preliminaries

Common to all problems treated in this paper is the input of a set of $m$ strings $S = \{S_1, \ldots, S_m\}$. For simplicity, we assume that all strings have the same length $n$, and that all characters are drawn from an alphabet $\Sigma$ of size $\sigma = |\Sigma|$. Hence, $|S_x| = n$ denotes the length of each input string and $S_x[i] \in \Sigma$ for all $i \in [1..n]$ and $x \in [1..m]$. Except for MCSP, the output is a string $T$ that is object to an optimization argument with respect to the input strings (and, additionally for CSS, with respect to an integer parameter specifying the length of $T$).

### Encoding Annotations.

Beginning with the next section, we state rules and constraints with numbered equations, and add to each equation, in square brackets, the number of generated clauses and the size of each such clause. For instance, the equation

$$[O(n), O(1)] \quad \forall i \in [1..n]: p_i \implies p_{i+1}$$

defines $n$ clauses, each of the form $(\neg p_i \lor p_{i+1})$, so its complexity is $[O(n), O(1)]$.

### Experiments.

We implemented our MAX-SAT-formulations in the ASP language, and used the solver clingo \cite{26, 27}\footnote{https://github.com/potassco/clingo} for evaluation. We compare the results with brute-force approaches written in the python language on randomly generated data. Our filenames are formatted like s03m04n005i1 to denote that the alphabet size is $\sigma = 3$, the number of strings is $m = 4$, the length of each string is $n = 5$, and this file is the $i = 1$-st sample of a batch of files with the same characteristics ($\sigma, m$ and $n$). For MCSP, we have file formats like 2s02n008i2.txt where the prefix 2 denotes that $m = 2$ is fixed. For the MCSP files, we assume that the two strings given have the same Parikh vector. Our implementations and datasets are available at https://github.com/koeppl/aspstring. For the evaluation, all experiments ran single-threaded on a machine with Intel Core i3–9100 CPU and Debian 11.
We use the known fact that we have to select, for the \(i\)-th character of the output \(T\), a character appearing at the \(i\)-th position of one of the input strings.

**Lemma 1 (\cite{37, Lemma 2}).** For each \(i \in [1..n]\), \(T[i] = S_x[i]\) for an \(x \in [1..m]\).

Let us define \(\Sigma_i := \{S_1[i], \ldots, S_m[i]\}\) to be the set of characters appearing at text position \(i\) of all input strings. Then \(\sigma_i := |\Sigma_i| \leq \min(m, \sigma)\), and \(\sigma_i\) can be much less than \(m\) or \(\sigma\) if the number of distinct characters is small. We can express the alphabets per position \(\Sigma_i\) by a Boolean matrix \(M[1..n][1..\sigma]\) with \(M[i][c] = 1\) if \(c \in \Sigma_i\).

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4 Alternative names are, among others, **Minimum Radius, Center String or Consensus String** problem.
Further, we define the variables $T_{i,c} = 1$ to encode that $T[i] = c$, for $i \in [1..n], c \in \{S_1[i], \ldots, S_m[i]\}$. To state that $T[i] = S_x[i]$, we want that, for a fixed position $i \in [1..n]$, only one $T_{i,c}$ is set:

$$[O(n), O(\min(m, \sigma))] \forall i \in [1..n] : \sum_{c \in \Sigma} T_{i,c} = 1 \quad \text{(CSP1)}$$

Next, we define the cost variables $C_{i,x}$ for all $i \in [1..n]$ and $x \in [1..m]$ with $C_{i,x}$ being set if $T[i] \neq S_x[i]$. Thus the Hamming distance between $T$ and $S_x$ is $\text{dist}_{\text{ham}}(T, S_x) = \sum_{i \in [1..n]} C_{i,x}$. Therefore:

$$[O(nm\sigma), O(1)] \forall i \in [1..n], c \in \Sigma, x \in [1..m] : T_{i,c} \land S_x[i] \neq c \implies C_{i,x} \quad \text{(CSP2)}$$

A statement for setting $C_{i,x}$ to false is not needed as the optimizer will try to do so if it does not violate (CSP2). This is achieved by the following objective:

$$[O(1), O(mn)] \text{minimize } \max_{x \in [1..m]} \sum_{i \in [1..n]} C_{i,x} \quad \text{(CSP3)}$$

**Complexities.** We have $O(n\sigma)$ selectable variables ($T_{i,c}$), $O(nm)$ helper variables ($C_{i,x}$), $O(nm\sigma)$ clauses (CSP2). The largest clause contains $O(mn)$ variables (CSP3).

**Implementation.** Our implementation in ASP is given in Listing 1. In all listings, the percent sign % introduces a comment until the end of the line, which we use to refer to the MAX-SAT equation that is represented by the respective line of code. Red curly arrows symbolize line breaks. If not otherwise stated, in all code listings onwards, we assume that the input is of the form $s(X, I, C)$, denoting that $S_x[I] = C \in \Sigma$. We use the helper variables $\text{mat}(X, I)$ to denote the existence of $S_x[I]$. For encoding (CSP3) in ASP, we additionally define the helper variables $\text{cost}$ and $\text{mcost}$ encoding $\sum_{i \in [1..n]} C_{i,x}$ and $\max_{x \in [1..m]} \text{cost}(x)$, respectively. The #show directives at the end define the variables the solver has to output. The evaluation for our implementation is deferred until we have introduced the css problem, which we jointly evaluate in Sect. 4.2.

```
Listing 1 ASP for csp (Sect. 3).
mat(X, I) :- s(X, I, C).
1 \{ t(I, C) : s(_, I, C) \} 1 :- mat(_, I). \%(CSP1)
c(X, I) :- t(I, C), s(X, I, A), C != A. \%(CSP2)
cost(X, I) :- C = \#sum \{1, l : c(X, I)\}, mat(X, _). \%(CSP3)
mcost(M) :- M = \#max \{C : cost(_, C)\}.
#minimize \{M : mcost(M)\}.
#show t/2. #show mcost/1. #show cost/2.
```

### 4 Closest Substring (CSS)

For the css problem, we additionally require a parameter $\lambda$ as input to specify the length of the output string $T$. css asks for the string $T$ with $|T| = \lambda$ such that $\max_{x \in [1..m]} \text{dist}_{\lambda}(S_x, T)$ is minimal, where $\text{dist}_{\lambda}(S_x, T) := \min_{i \in [1..n-\lambda+1]} \text{dist}_{\text{ham}}(S_x[i..i+\lambda-1], T)$ is the number of mismatches we need to be able to detect $T$ via approximate pattern matching in $S_x$ with $\text{dist}_{\lambda}(S_x, T)$ mismatches. An example is shown in Fig. 2.
Figure 2 Example for css (Sect. 4) with $n = 13$ and query length $\lambda = 4$. The input is shown on the left figure. We can observe in the right figure that $T = \text{snes}$ is the css having one mismatch with each of the input strings in the Hamming distance by horizontally shifting the input strings.

**Related Work.** The decision problem for $\delta$ mismatches is also called $\delta$-MISMATCH problem. Gramm et al. [32, Theorem 2] solved the decision problem in $O(m\lambda + (n - \lambda)m^{\delta + 1})$ time. Marx et al. [46] showed that css can be solved in $O(\lambda \theta)O(n(nm\delta + 1)\theta)O(\log\delta)\theta)$ or $O((\sigma\delta)O(n(m)O(log\log m))$ time. A survey on further results can be found in [31]. With respect to other optimization approaches, we are aware of a genetic algorithm [47].

4.1 MAX-SAT encoding

Following [32, Section 3.3], we reduce css to CSP by selecting shifts $d_x \in [0..n - \lambda]$ of each input string $S_x$ such that the CSP of $\{S_x[1 + d_1, \ldots, \lambda + d_1], \ldots, S_m[1 + d_m, \ldots, \lambda + d_m]\}$ is a solution of css if we take the minimum distance over all shifts $d_x$.

In what follows, we represent the shifts by a matrix of selectable Boolean variables of size $O(m(n - \lambda))$. We redefine the alphabet for the $i$-th character to be $\Sigma_i := \{S_x[i + d_i], \ldots, S_m[i + d_m]\}$. We define the variables $T_{i,c}$ and $C_{i,x}$ as before. We copy (CSP1) as it is since it only states from which string $S_x$ we select the $i$-th character of $T$, except that we have $O(\lambda)$ instead of $O(n)$ clauses since $|T| = \lambda$. The major difference is that for checking equality, we must add the offsets and obtain the following modification of (CSP2):

$$[O(\lambda m\sigma), O(1)] \forall i \in [1..\lambda], c \in \Sigma_i, x \in [1..m] : T_{i,c} \land S_x[i + d_x] \neq c \implies C_{i,x} \quad \text{(CSS2)}$$

The additional $n$-term in the complexity stems from the fact that the offsets $d_x$ are represented as a two-dimensional binary array. The other equations as well as the objective are kept in the same way.

**Complexities.** We have $O(\lambda \sigma + m(n - \lambda))$ selectable variables ($T_{i,c}$ and $d_x$), $O(\lambda m)$ helper variables ($C_{i,x}$), $O(\lambda m n \sigma)$ clauses. The largest clause has size $O(\lambda m)$. Our implementation in ASP is given in Listing 2, where we expect an additional input of the form `#const lambda=\lambda`. for the requested substring length $\lambda$.

**Listing 2 ASP for css (Sect. 4).**

```
mat(X,I) :- s(X,I,\_).  
1 {d(X,D) : D = 0..n-lambda} 1 :- mat(X,0).  
sigma(I,C) :- s(X,J,C), d(X,D), J-D >= 0, I = J-D.  
1 {t(I,C) : sigma(I,C)} 1 :- mat(_,I), I < lambda.  
% (CSP1)

1 c(X,I) :- t(I,C), s(X,J,A), d(X,D), I+D = = J, I < lambda, A != C.  
% (CSS2)

cost(X,C) :- C = #sum {1,I : c(X,I)}, mat(X,\_).  
% (CSP2)

mcost(M) :- M = #max {C : cost(_,C)}.  
% (CSP3)

#minimize {M : mcost(M)}.  
#show t/2.  #show mcost/1.  #show cost/2.
```
### Encoding Hard String Problems with Answer Set Programming

Table 1 Evaluation for the CLOSEST STRING PROBLEM (csp) for \( \lambda = 0 \) and CLOSEST SUBSTRING PROBLEM (css) for \( \lambda > 0 \). The column \( \text{dist} \) shows the maximum Hamming distance of the reported string to all input strings. The column \( \text{rules} \) is the number of created SAT rules, \( \text{vars} \) is the number of variables, and \( \text{choices} \) is the number of choices or configurations the solver or brute-force algorithm tries. Reported times are in seconds (s).

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<th>rules</th>
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</table>

#### 4.2 Evaluation of csp and css

Although there are efficient heuristics like choosing a majority string [8], we compared our ASP encoding for csp to a basic brute-force algorithm that enumerates all possible assignments for the characters of the closest substring. The number of possible configurations for \( T \) is \( e_{SS} := \prod_{i=1}^{n} \sigma_i \in O(\min(\sigma^n), \lambda \sigma^n) \) dependent on the shape of the strings in \( S \). A brute-force algorithm trying each configuration spends \( O(e_{SS} m^n) \) time on computing the Hamming distances of the resulting string \( T \) with all strings of \( S \).

This algorithm can be easily adopted for css. For that, we consider all possible offsets of the input strings like in the ASP encoding. Hence, the number of configurations is the number of configurations for the csp instance, multiplied by \((n - \lambda)^m\) for each possible offset value. If \( \lambda \) is small, then it suffices to compute all configurations of \( T \), which are \( \sigma^\lambda \) many, and compute the Hamming distances in \( O(\lambda m) \) time for each such configuration. We implemented the former brute-force approach, whose time complexity grows exponentially with all parameters \( \sigma, n, \) and \( m \), for randomly generated strings. We can observe this case in Table 1, where the ASP implementation outperforms the brute-force approach.
Table 2: Evaluation of the Closest String problem (scp) on datasets provided by Torres and Hoshino [54]. The column distance is the maximal Hamming distance of the output to any of the input strings.

<table>
<thead>
<tr>
<th>file</th>
<th>distance</th>
<th>rules</th>
<th>vars</th>
<th>choices</th>
<th>time [s]</th>
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<td>8</td>
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<td>1334</td>
<td>18,095</td>
<td>0.373</td>
</tr>
<tr>
<td>rand-20-50-50-10-1</td>
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<td>3574</td>
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<td>1.255</td>
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<tr>
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<td>19.168</td>
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<tr>
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<td>24,654</td>
<td>6162</td>
<td>39,265,311</td>
<td>7009.909</td>
</tr>
</tbody>
</table>

In Table 2, we depict the results of a larger evaluation on the datasets provided in [54][5], which are also used in [16, 43]. We kept their file naming, which is the format `rand-σ-m2-n2-i`, where `i` is an iteration counter to have multiple files with the same characteristics (`m`, `n`, and `σ`). The prefix `rand` can be followed by a zero. We observe that larger distances correlate with the number of choices, affecting the overall running time. Even for large inputs with short distances like the dataset `rand-4-150-150-5-1`, the running time is short.

Figure 3 Example for LCS (Sect. 5) with \( n = 13 \). The input is shown on the left figure. In the right figure, we highlighted the subsequences matching \( T = \text{sleepless} \) by surrounding the respective characters with boxes in each input string. Here, \( T = \text{sleepless} \) is the LCS of all input strings.

5 Longest Common Subsequence (LCS)

The LCS problem asks for the longest string \( T \) such that \( T \) is a subsequence of \( S_x \) for every \( x \in [1..m] \). See Fig. 3 for an example.

Existence. A solution exists if all strings share at least one common character in the alphabet.

Related Work. Maier [45] showed that LCS is NP-hard for \( \sigma \geq 2 \), and the same holds for SCS with \( \sigma \geq 5 \). Later, Blin et al. [5] gave a proof that LCS stays NP-hard even if the input strings are well-compressible with the run-length encoding. For exact algorithms, we can extend the classic dynamic programming (DP) algorithm of Wagner and Fischer [56] to \( m \) strings, which then takes \( O(n^m) \) time. Irving and Fraser [36] gave two algorithms running in \( O(mn(n-\ell)^{m-1}) \) or \( O(m\ell(n-\ell)^{m-1}+m\sigma n) \) time, where \( \ell \) is the length of the output. This result implies that LCS is FPT in \( m \) and \( n-\ell \). Bulteau et al. [10] improved the result of [36] with an algorithm running in \( O((n-\ell+1)^{n-\ell+1}mn) \) time, which is an FPT in the number of deletions \( n-\ell \). Finally, there is a genetic algorithm [34] and an ant colony optimization algorithm [50].

5.1 MAX-SAT encoding

Our idea is to select a subsequence \( T_x \) for each input string \( S_x \) and maximize the length of \( T_x \) under the constraint that all \( T_x \)’s have to be equal. The subsequence \( T_x \) of \( S_x \) is given by a sequence of indices \( i_1 < \ldots < i_{|T_x|} \) such that \( S_x[i_1] \cdots S_x[i_{|T_x|}] = T_x \). We can encode the subsequences \( T_x \) by the selectable variables \( C_{x,\ell,i} \) encoding whether \( T_x[\ell] = S_x[i] \), for each \( x \in [1..m], \ell \in [1..n] \). We make use of \( C_{x,\ell,i} \) as follows. First, for each \( T_x[\ell] \), we define the range for the selectable variables \( C_{x,\ell,i} \).

\[ \forall x \in [1..m], \ell \in [1..n] : \sum_{i \in \ell..n} C_{x,\ell,i} \geq 0 \quad \text{(LCS1)} \]

\[ \text{Logically, we would expect in (LCS1) a "\leq 1" instead of a "\geq 0". However, the former suffices together with the following constraints and is cheaper than "\leq 1".} \]

\[ [O(nm), O(n)] \]
If we have selected $T_x[\ell]$ to be $S_x[i]$, then $T_x[\ell-1]$ must be a character chosen in $S_x[1..i-1]$:

$$[\mathcal{O}(n^2 m), \mathcal{O}(n)] \quad \forall x \in [1..m], \ell \in [2..n], i \in [\ell..n] :$$

$$C_{x,\ell,i} \implies \sum_{j \in [1..i-1]} C_{x,\ell-1,j} = 1 \quad (\text{LCS2})$$

Next, we define the helper variables $V_{x,\ell}$ encoding whether $T_x$ has a length of at least $\ell$, for each $x \in [1..m], \ell \in [1..n]$. If we have selected a character for $T_x[\ell]$ via $C_{x,\ell,i}$, then we set $V_{x,\ell}$ to true to specify that $T_x$ has a length of at least $\ell$.

$$[\mathcal{O}(nm), \mathcal{O}(n)] \quad \forall x \in [1..m], \ell \in [1..n] : \bigvee_{i \in [1..n]} C_{x,\ell,i} \implies V_{x,\ell} \quad (\text{LCS3})$$

We now restrict all $T_x$'s to be of equal length, which we do in a Round-Robin encoding:

$$[\mathcal{O}(nm), \mathcal{O}(1)] \quad \forall x \in [1..m], \ell \in [1..n] : V_{x,\ell} \implies V_{(x+1) \bmod n, \ell} \quad (\text{LCS4})$$

Here, $\bmod n : \{1, 2, \ldots\} \rightarrow [1..n]$ is the modulo operation with $n \bmod n = n$ and $(n + 1) \bmod n = 1$. To achieve that all $T_x$ store the same characters, we use the following constraint.

$$[\mathcal{O}(n^3 m), \mathcal{O}(1)] \quad \forall x \in [1..m], \ell \in [1..n], i, j \in [1..n] :$$

$$C_{x,\ell,i} \land C_{(x+1) \bmod m, \ell,j} \implies S_x[i] = S_{(x+1) \bmod m}[j] \quad (\text{LCS5})$$

Finally, we enforce that we need to select a position for $T_x[\ell]$ if $V_{x,\ell}$ is set:

$$[\mathcal{O}(nm), \mathcal{O}(n)] \quad \forall x \in [1..m], \ell \in [1..n] : V_{x,\ell} \implies \bigvee_{i \in [\ell..n]} C_{x,\ell,i} \quad (\text{LCS6})$$

Alternatively to (LCS5) and (LCS6), we can state that the next subsequence must select one of the text positions $j$ for $T_{x+1}[\ell]$ with $S_{x+1}[j] = S_x[i]$.

$$[\mathcal{O}(n^2 m), \mathcal{O}(n)] \quad \forall x \in [1..m], \ell \in [1..n], i \in [1..n] :$$

$$C_{x,\ell,i} \implies \sum_{j : S_x[i] = S_{(x+1) \bmod n}[j]} C_{(x+1) \bmod m, \ell,j} = 1 \quad (\text{LCS5}')$$

Finally, we formulate our optimization problem as

$$[\mathcal{O}(1), \mathcal{O}(n)] \quad \text{maximize} \sum_{\ell \in [1..n]} V_{1,\ell} \quad (\text{LCS7})$$

**Complexities.** Our implementation in ASP is given in Listing 3. We have $\mathcal{O}(nm^2)$ selectable variables ($C_{x,\ell,i}$), $\mathcal{O}(nm)$ helper variables ($V_{x,\ell}$), and $\mathcal{O}(n^2 m)$ clauses (LCS5'). The largest clause has $\mathcal{O}(n)$ variables. An improvement for short LCS solutions could be to encode the existence problem for a fixed length $\lambda$ in ASP such that we have $\mathcal{O}(m\lambda)$ selectable variables for encoding $T_x$, and call this encoding while varying $\lambda$ to find the largest value for $\lambda$ admitting a solution.
Table 3 Evaluation of the LONGEST COMMON SUBSEQUENCE problem (lcs).

<table>
<thead>
<tr>
<th>file</th>
<th>length</th>
<th>rules</th>
<th>vars</th>
<th>choices</th>
<th>time [s]</th>
<th>choices</th>
<th>time [s]</th>
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<td>8388608</td>
<td>43.35</td>
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</tr>
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<td>9640</td>
<td>0.47</td>
<td>4194304</td>
<td>15.94</td>
</tr>
</tbody>
</table>

Listing 3 ASP for lcs (Sect. 5).

```prolog
mat(X,I) :- s(X,I,\_).
0 {c(X,L,I) : mat(X,I), I >= L} :- mat(X,L).
% (LCS1)
1 {c(X,L,J) : J < I, mat(X,J)} :- c(X,L+1,I), mat(X,L), mat(X,L+1).
% (LCS2)
v(X,L) :- c(X,L,I), mat(X,I), mat(X,L).
% (LCS3)
v(X+1,L) :- v(X,L), mat(X,L), mat(X+1,L).
% (LCS4)
v(0,L) :- v(m-1,L).
% (LCS5)
:- c(X+1,L,J), c(X,L,I), s(X,I,D), not s(X+1,J,D).
% (LCS5)
:- c(0,L,J), c(m-1,L,I), s(m-1,I,D), not s(0,J,D).
1 {c(X,L,I) : mat(X,I), I >= L} :- v(X,L).
% (LCS6)
#maximize {1,L : v(0,L)}.  % (LCS7)
#show c/3.
```

5.2 Evaluation

A DP approach would need $O(n^m)$ time (cf. [15, Chapter IV, Section 15.4] for a textbook reference). Here, we stick to a trivial approach that tries all distinct subsequences of the first string $S_1$, and for each such subsequence we check whether it is a subsequence of all other input strings. The number of these subsequences is at most $2^n - 1$. If we select these subsequences with respect to their lengths, starting with the longest possible one, we can terminate whenever the selected subsequence is found in all other strings. In the worst
case, the time complexity of this approach grows exponentially in \( n \), but only linearly in \( m \), independent of the alphabet size. We therefore restrict our evaluation in Table 3 to scaling \( n \) while keeping the other parameters unchanged. Like in Sect. 4.2, the ASP implementation outperforms the brute-force approach. However, a DP implementation might outperform the ASP implementation by re-using memoized results.

### 6 Minimum Common String Partition (MCSP)

For the special case of \( m = 2 \) input strings \( S_1 \) and \( S_2 \), the MCSP problem, introduced by Goldstein et al. [29] and Swenson et al. [52], asks, for a given \( z \in [1..n] \), a factorization of \( S_1 \) into \( S_x = F_1 \cdots F_z \) and a permutation \( \pi \) of \([1..z]\) such that \( F_{\pi(1)} \cdots F_{\pi(z)} = S_2 \). The optimization problem is to find the smallest \( z \) for which a solution exists. We give an example in Fig. 4.

**Existence.** A sufficient condition for whether a solution for any \( z \in [1..n] \) exists is to check that the Parikh vectors of \( S_1 \) and \( S_2 \) are the same, such that at least a permutation on \([1..n]\) exist to rearrange the characters of \( S_1 \) to form \( S_2 \).

**Related Work.** While introducing MCSP, Goldstein et al. [29] also showed that it is NP-hard. Bulteau and Komusiewicz [11] showed that MCSP is FPT in \( z \). For constant alphabets (\( \sigma = O(1) \)), Cygan et al. [17] presented an exact algorithm running in \( 2^{O(n \lg n / \lg \lg n)} \) time. Recently, Chromý and Sinnl [14] studied a DP algorithm. It is known that MCSP can be tackled by probabilistic tree searches [7], ILP formulations [6, 23], and an ant colony optimization algorithm [22].

### 6.1 MAX-SAT encoding

We adapt the MAX-SAT encoding of Bannai et al. [1] for the shortest bidirectional macro scheme problem [51]. To this end, we define the sets \( M_i := \{ j \in [1..n] \mid S_1[i] = S_2[j] \} \subset [1..n] \) specifying the positions in \( S_2 \) that match with \( S_1[i] \). In what follows, we make use of the following selectable Boolean variables:

- \( p_i \) or \( q_i \) encode if \( S_1[i] \) or \( S_2[i] \) is the start of a factor, respectively, for \( i \in [1..n] \).
- \( \text{ref}_{i \rightarrow j} \) encodes whether position \( i \) of \( S_1 \) references position \( j \) of \( S_2 \), and vice versa, for \( i \in [1..n] \) and \( j \in \mathcal{M}_i \).
We have $O(n^2)$ Boolean variables, which we use as follows. On the one hand, each position in $S_1$ has exactly one reference:

$$\forall i \in [1..n]: \sum_{j \in \mathcal{M}_i} \text{ref}_{i \rightarrow j} = 1$$

(MCSP1)

On the other hand, each position in $S_2$ has exactly one reference:

$$\forall j \in [1..n]: \sum_{i \in [1..n]} \text{ref}_{i \rightarrow j} = 1$$

(MCSP2)

In what follows, we add implications for the factor starting positions that are due to how we set the references. First, a factor starts always at the first text position, so $p_1$ and $q_1$ are always true. If $S_1[i]$ references $S_2[i]$ and $i$ is a factor starting position of $S_1$, so is $j$ for $S_2$.

$$\forall i \in [1..n], j \in \mathcal{M}_i: p_i \land \text{ref}_{i \rightarrow j} \implies q_j$$

(MCSP3)

Next, if $S_1[i]$ references $S_2[i]$ and $j$ is a factor starting position of $S_2$, so is $i$ for $S_1$. We only have to check that condition for $q_1$ since all other constraints setting $p_i$ and constraint (MCSP3) then implies that $q_j$ has to be set.

$$\forall i \in [1..n]: q_1 \land \text{ref}_{i \rightarrow 1} \implies p_i$$

(MCSP4)

Another condition is that if the previous text positions have mismatching characters, we cannot extend the factor to the left.

$$\forall i \in [1..n], j \in \mathcal{M}_i \text{ with } S_1[i-1] \neq S_2[j-1]: \text{ref}_{i \rightarrow j} \implies p_i$$

(MCSP5)

Even if the previous characters match, when the reference of the previous text positions is different, we need to make a factor starting position:

$$\forall i \in [2..n], j \in \mathcal{M}_i \text{ such that } j > 1 \text{ and } S_2[i-1] = S_2[j-1]: 
\neg \text{ref}_{i-1 \rightarrow j-1} \land \text{ref}_{i \rightarrow j} \implies p_i$$

(MCSP6)

Finally, we minimize

$$\sum_{i \in [1..n]} p_i$$

(MCSP7)

**Complexities.** We have $O(n^2)$ selectable variables, and $O(n^2)$ clauses (MCSP3). The largest clause has $O(n)$ variables (MCSP2). Our implementation in ASP is given in Listing 4. Note that we start counting at zero, so $p(0)$ is equivalent to setting $p_1$ to true. Instead of mat we use the helper variables spos and tpos denoting the existence of $S_1[i]$ and $S_2[i]$, respectively.

**Listing 4** ASP for MCSP (Sect. 6).

```plaintext
spos(0) :- s(0, I, _).
tspos(J) :- s(1, J, _).
p(0). q(0).
arc(I, J) :- s(0, I, C), s(1, J, C).
1 {ref(I, J) : arc(I, J)} 1 :- spos(I). % (MCSP1)
1 {ref(I, J) : arc(I, J)} 1 :- tpos(J). % (MCSP2)
q(J) :- p(I), ref(I, J). % (MCSP3)
p(I) :- q(I), ref(I, I). % (MCSP4)
p(I) :- ref(I, J), s(0, I-1, C), s(1, J-1, D), C != D. % (MCSP5)
p(I) :- not ref(I-1, J-1), ref(I, J). % (MCSP6)
#minimize {1, X : p(X)}.
#show ref/2. #show p/1. #show q/1.
```
Table 4 Evaluation of the Minimum Common String Partition problem (mcsp). Note that the time for the ASP solution is in milliseconds. The column $z$ denotes the number of factors of the returned partition.

<table>
<thead>
<tr>
<th>file</th>
<th>$z$</th>
<th>rules</th>
<th>vars</th>
<th>choices</th>
<th>time [ms]</th>
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<td>25</td>
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<td>61</td>
<td>2.0</td>
<td>986409</td>
<td>6.31</td>
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<td>2s02n010i0</td>
<td>4</td>
<td>510</td>
<td>213</td>
<td>108</td>
<td>2.0</td>
<td>9864100</td>
<td>70.92</td>
</tr>
<tr>
<td>2s03n010i0</td>
<td>6</td>
<td>484</td>
<td>147</td>
<td>37</td>
<td>1.0</td>
<td>9864100</td>
<td>71.04</td>
</tr>
<tr>
<td>2s03n010i2</td>
<td>4</td>
<td>584</td>
<td>164</td>
<td>47</td>
<td>2.0</td>
<td>9864100</td>
<td>71.38</td>
</tr>
<tr>
<td>2s02n010i2</td>
<td>4</td>
<td>637</td>
<td>189</td>
<td>77</td>
<td>2.0</td>
<td>9864100</td>
<td>73.78</td>
</tr>
<tr>
<td>2s02n010i1</td>
<td>3</td>
<td>639</td>
<td>187</td>
<td>103</td>
<td>2.0</td>
<td>9864100</td>
<td>74.28</td>
</tr>
</tbody>
</table>

Table 5 Evaluation of the Minimum Common String Partition problem (mcsp) on prefixes of the SARS-CoV-2 dataset.

<table>
<thead>
<tr>
<th>length</th>
<th>$z$</th>
<th>rules</th>
<th>vars</th>
<th>choices</th>
<th>time [s]</th>
</tr>
</thead>
<tbody>
<tr>
<td>10</td>
<td>4</td>
<td>447</td>
<td>146</td>
<td>34</td>
<td>0.001</td>
</tr>
<tr>
<td>20</td>
<td>12</td>
<td>1273</td>
<td>445</td>
<td>269</td>
<td>0.003</td>
</tr>
<tr>
<td>30</td>
<td>14</td>
<td>2282</td>
<td>911</td>
<td>1951</td>
<td>0.017</td>
</tr>
<tr>
<td>40</td>
<td>16</td>
<td>3720</td>
<td>1685</td>
<td>4683</td>
<td>0.047</td>
</tr>
<tr>
<td>50</td>
<td>21</td>
<td>5468</td>
<td>2442</td>
<td>2050092</td>
<td>18.609</td>
</tr>
<tr>
<td>60</td>
<td>24</td>
<td>7451</td>
<td>3422</td>
<td>6866999</td>
<td>80.256</td>
</tr>
</tbody>
</table>

6.2 Evaluation

Without leveraging the actual contents of the characters like in our SAT formulation, a naive way is to factorize both strings $S_1$ and $S_2$ with factors of the same lengths, and check whether there exists a permutation such that we can match factors of $S_1$ with factors of $S_2$. To this end, we iterate over the size $z$ of the partition from 1 to $n$. For each $z \in [1..n]$, we partition $S_1$ into $z$ factors $S_1 = F_1 \cdots F_z$. There are $\binom{n}{z}$ such ways to partition $S_1$. For each permutation $\pi_z$ on $[1..z]$, we define the factorization $G_1 \cdots G_z = S_2$ with $|G_x| = |F_{\pi(x)}|$ for all $x \in [1..z]$. If $G_x = F_{\pi(x)}$ for all $x \in [1..z]$, then we have found a solution, and terminate. The number of configurations is $\sum_{z=1}^{n} \binom{n}{z} z!$, and each check takes $O(n)$ time. Like the brute-force approach for LCS (Sect. 5.2), this approach has an exponential dependency on the text length $n$. In Table 4, we observe that specifying the choices for the references for each position individually (as we do in our ASP encoding) reduces the number of choices significantly when compared to the choices the brute-force algorithm processes.

Since our ASP encoding for mcsp seems quite efficient, we subsequently performed a benchmark on real data. In detail, we conducted an experiment by scaling the prefix length of a given input sequence, and report results in Table 5. For that, we used the SARS-CoV-2
The scs problem asks for the shortest string $T$ such that $S_x$ is a substring of $T$, for all $x \in [1..m]$. Figure 5 shows an example.

**Existence.** A trivial common superstring is the concatenation $S_1 \cdots S_m$. Permuting the strings and removing overlapping parts lead to the solution [25].

**Related Work.** Gallant et al. [25] showed that scs is NP-hard for $n \geq 3$ with respect to the number of strings $m$ and unbounded alphabet size, but can be solved in linear time if $n \leq 2$. For binary alphabet $\sigma = 2$, they showed that the problem is still NP-hard for $n = \Omega(\log(nm))$.

### 7.1 Reduction to Hamiltonian Path

We follow the idea of Tarhio and Ukkonen [53] by reducing scs to the search of the Hamiltonian path maximizing the weights of the selected edges. The ASP encoding of finding a Hamiltonian cycle in an unweighted graph has already been studied in [42, 41]. We build on one of their approaches and extend it by maximizing the weights while omitting the weight of one edge to turn the cycle into a Hamiltonian path\(^8\). An overlap graph $(S, A, w)$ is a weighted

---

\(^7\) [https://github.com/cfarkas/SARS-CoV-2-freebayes](https://github.com/cfarkas/SARS-CoV-2-freebayes)

\(^8\) We make a distinction between Hamiltonian path and Hamiltonian cycle in the sense that the cycle visits exactly one node twice.
directed graph, having the input strings $S$ as nodes and the arcs $A := \{(S_x, S_y) : x \neq y\}$. The weights are defined by a weight function $w : A \to [0..n]$ with $w(S_x, S_y) := \max\{|U| : U \text{ is suffix of } S_x \text{ and prefix of } S_y\}$. Hence, $w(S_x, S_y)$ is the number of overlapping characters, which we can omit if we want to build the superstring of $S_x$ and $S_y$ that starts with $S_x$.

With respect to the overlap graph, a path is a sequence of strings, and a Hamiltonian path in the overlap graph is a path that visits each node exactly once, i.e., a permutation $\pi$ of the list $\{S_1, \ldots, S_m\}$. Our goal is to find a permutation that maximizes $\sum_{x=1}^{m-1} w(S_{\pi(x)}, S_{\pi(x+1)})$, i.e., to find the Hamiltonian path whose arcs have maximal weights in sum.

7.2 MAX-SAT encoding

We define the following $O(m^2)$ Boolean variables:

- $cycle_{x,y}$ encoding whether we have the arc $(S_x, S_y)$ in our Hamiltonian cycle, for $x, y \in [1..m]$;
- $reach_{x,y}$ encoding whether we can reach $S_y$ from $S_x$ by following the transitive closure of $cycle$, for $x, y \in [1..m]$;
- $start_x$ encoding whether our superstring starts with $S_x$, for $x \in [1..m]$.

First, we select arcs from the overlap graph for $cycle_{x,y}$. To this end, for each string $S_x$, we select exactly one out-going arc and one in-coming arc:

$$[O(m), O(m)] \forall x \in [1..m] : \sum_{y=1}^{m} cycle_{x,y} = 1 \text{ and } \forall y \in [1..m] : \sum_{x=1}^{m} cycle_{x,y} = 1 \quad (SCS1)$$

The transitive closure of $cycle$ can be encoded as follows. First we initialize $reach$ by the direct connections due to $cycle$.

$$[O(m^2), O(1)] \forall x, y \in [1..m], x \neq y : cycle_{x,y} \implies reach_{x,y} \quad (SCS2)$$

Next, if we can reach $y$ from $x$, and there is an arc $(y, z)$, then we can reach $z$ from $x$:

$$[O(m^3), O(1)] \forall x, y, z \in [1..m], x \neq y \neq z : reach_{x,y} \wedge cycle_{y,z} \implies reach_{x,z} \quad (SCS3)$$

To make the path selected by $cycle_{x,y}$ an Hamiltonian path, we want that all strings are connected via $reach$:

$$[O(m^2), O(1)] \forall x, y \in [1..m], x \neq y : reach_{x,y} = 1 \quad (SCS4)$$

For the Hamiltonian path it is left to select a designated start string$^9$.

$$[O(1), O(m)] \sum_{y=1}^{m} start_y = 1 \quad (SCS5)$$

Finally, our objective is to maximize the weights on the path starting from $start_x$ of length $m$:

$$[O(1), O(m^2)] \text{ maximize } \sum_{x,y \in [1..m]} w(x, y) \quad (SCS6)$$

$^9$ It actually suffices to check in (SCS4) that all strings can be reached from this start string, but doing so had a negative impact on the overall running time in the experiments.
Table 6 Evaluation of the Shortest Common Superstring problem (scs). $|T|$ is the length of the SCS.

|     | $|T|$  | rules | vars  | choices | time [s] | choices | time [s] |
|-----|-------|-------|-------|---------|----------|---------|----------|
| s02m10n008i0 | 42    | 2090  | 1416  | 198756  | 3.58     |         |          |
| s02m10n008i1 | 33    | 2206  | 1465  | 1854941 | 40.73    | 10240   | 0.02     |
| s02m10n008i2 | 39    | 2200  | 1464  | 1401686 | 29.49    | 10240   | 0.02     |
| s02m11n008i0 | 49    | 2639  | 1825  | 2150681 | 44.96    | 22528   | 0.03     |
| s02m11n008i1 | 35    | 2699  | 1861  | 6652411 | 154.48   | 22528   | 0.03     |
| s02m11n008i2 | 50    | 2611  | 1817  | 6980725 | 136.00   | 22528   | 0.02     |

Complexities. We have $O(m^2)$ selectable variables and $O(m^3)$ clauses (SCS3). The largest clause has $O(m^2)$ variables (SCS6). Our implementation in ASP is given in Listing 5. We expect an input of the form $w(X, Y, C)$ encoding the weight $w(X, Y) = C$. The helper variables node and gain define the nodes of the overlap graph and the value of the optimization argument in (SCS6), respectively.

Listing 5 ASP for scs (Sect. 7).
```
node(X) :- w(X,_,_).
1 {cycle(X,Y) : w(X,Y,_)} 1 :- node(X).  \%(SCS1)
1 {cycle(X,Y) : w(X,Y,_)} 1 :- node(Y).
reach(X,Y) :- cycle(X,Y).  \%(SCS2)
reach(X,Z) :- reach(X,Y), cycle(Y,Z).  \%(SCS3)
:- not reach(X,Y), node(X), node(Y).  \%(SCS4)
1 {start(X) : node(X)} 1.  \%(SCS5)
gain(D) :- D = #sum {C,X : cycle(X,Y), w(X,Y,C), not start(Y)}.  \%(SCS6)
#maximize {D : gain(D)}.
#show cycle/2.  #show start/1.
```

7.3 Evaluation

The overlap graph can be computed in $O(nm + m^2)$ time [33]. Given the overlap graph, the easiest approach is to enumerate all $m!$ permutations, and compute the sum of the selected weights in $\Theta(m)$ time. The time bound can be improved by using a DP approach taking $O(m^2 2^m)$ time\(^{10}\). In the experiments of Table 6, we use this DP approach as our brute-force solution. We observe that it outperforms our ASP implementation on all instances. That is due to the fact that (a) our ASP encoding does not make use of more information than the DP approach, and that (b) the number of choices in our encoding for the Hamiltonian path is prohibitively large. As a matter of fact, efficient SAT and ASP encodings for Hamiltonian cycles are actively studied, cf. [58] for SAT and [4] for ASP.

\(^{10}\)https://leetcode.com/problems/find-the-shortest-superstring/solutions/194891/official-solution/
Table 7 Encoding complexities of the studied problems. Columns prob., #sel. vars, #h.vars, #clauses and max. cl. denote the problem name, the number of defined selectable variables, the number of helper variables, the number of clauses, and the maximum size a clause can have.

<table>
<thead>
<tr>
<th>prob.</th>
<th>#sel. vars</th>
<th>#h.vars</th>
<th>#clauses</th>
<th>max. cl.</th>
</tr>
</thead>
<tbody>
<tr>
<td>CSP</td>
<td>$O(n\sigma)$</td>
<td>$O(nm)$</td>
<td>$O(nm\sigma)$</td>
<td>$O(nm)$</td>
</tr>
<tr>
<td>CSS</td>
<td>$O(\lambda \sigma + (n - \lambda)m)$</td>
<td>$O(\lambda m)$</td>
<td>$O(nm\sigma \lambda)$</td>
<td>$O(\lambda m)$</td>
</tr>
<tr>
<td>LCS</td>
<td>$O(n^2 m)$</td>
<td>$O(mn)$</td>
<td>$O(n^2 m)$</td>
<td>$O(n)$</td>
</tr>
<tr>
<td>MCSP</td>
<td>$O(n^2)$</td>
<td>$O(1)$</td>
<td>$O(n^2)$</td>
<td>$O(n)$</td>
</tr>
<tr>
<td>SCS</td>
<td>$O(m^2)$</td>
<td>$O(1)$</td>
<td>$O(m^3)$</td>
<td>$O(m^2)$</td>
</tr>
</tbody>
</table>

8 Conclusion

We provided encodings in ASP for five prominent examples of NP-hard problems in the field of stringology. We summarized the complexities of the encodings in Table 7. We observed that, on the one hand, by leveraging characteristics of the input data such as for MCSP, our solution is far superior than simple brute-force approaches that omit those characteristics. On the other hand, for SCS, we observed that if the problem can be easily reduced to instances of problems like finding a Hamiltonian path, DP approaches are already efficient enough to find the answer faster than an ASP solver. It therefore depends on the nature of the problem we study for whether an application of an ASP solver makes sense. Nevertheless, the programming in ASP is highly expressive as can be seen by the short program codes in Listings 1–5, and therefore can be understood as a tool for rapid prototyping. Other advantages of ASP solvers like clingo are that they can work in parallel, report approximate solutions when reaching a given timeout, and enumerate all solutions, provided that the specified constraints do not exclude one of them. An evaluation of those features is left as future work since it would go beyond the scope of this paper.

References


Encoding Hard String Problems with Answer Set Programming


For small values of $\lambda$, the offsets can be quite large. Here, we present an alternative encoding without the offsets. The resulting encoding has fewer variables, but has more variables that are subject to the optimization argument. In what follows, we can encode $T[1..\lambda]$ by the Boolean variables $T_{i,c}$ specifying with $T_{i,c} = 1$ that $T[i] = c$:

$$[O(\lambda), O(\sigma)] \forall i \in [1..\lambda] : \sum_{c \in \Sigma} T_{i,c} = 1 \quad (CSS1')$$

We now let the costs encode the offsets by the variables $C_{i,x,o}$ being set if $S_x[o + i] \neq T[i]$.

$$[O(\lambda nm\sigma), O(1)] \forall i \in [1..\lambda], c \in \Sigma, x \in [1..m], o \in [1..n - \lambda] : T_{i,c} \land S_x[i + o] \neq c \implies C_{i,x,o} \quad (CSS2')$$
Table 8 Used Entities.

<table>
<thead>
<tr>
<th>entity</th>
<th>meaning</th>
</tr>
</thead>
<tbody>
<tr>
<td>Σ</td>
<td>alphabet</td>
</tr>
<tr>
<td>σ</td>
<td>alphabet size, (σ =</td>
</tr>
<tr>
<td>(S)</td>
<td>set of input strings ({S_1, \ldots, S_m})</td>
</tr>
<tr>
<td>(m)</td>
<td>size of (S), i.e., (m =</td>
</tr>
<tr>
<td>(n)</td>
<td>length of an input string</td>
</tr>
<tr>
<td>(S_x)</td>
<td>input string</td>
</tr>
<tr>
<td>(T)</td>
<td>string to output</td>
</tr>
<tr>
<td>(ℓ)</td>
<td>length for a subsequence</td>
</tr>
<tr>
<td>(δ)</td>
<td>distance of the output to all (S_x)</td>
</tr>
<tr>
<td>(i, j)</td>
<td>indices for text positions in an input string</td>
</tr>
<tr>
<td>(x, y)</td>
<td>indices for an input string</td>
</tr>
<tr>
<td>(c)</td>
<td>character in (Σ)</td>
</tr>
</tbody>
</table>

The objective function becomes

\[
\begin{align*}
[O(1), O(mn^2)] & \quad \text{minimize} \quad \max_{x \in [1..m]} \min_{o \in [1..n-\lambda]} \sum_{i \in [1..n]} C_{i,x,o} \\
& \quad \text{(CSS3')}
\end{align*}
\]
On the Complexity of Parameterized Local Search for the Maximum Parsimony Problem

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Abstract

Maximum Parsimony is the problem of computing a most parsimonious phylogenetic tree for a taxa set $X$ from character data for $X$. A common strategy to attack this notoriously hard problem is to perform a local search over the phylogenetic tree space. Here, one is given a phylogenetic tree $T$ and wants to find a more parsimonious tree in the neighborhood of $T$. We study the complexity of this problem when the neighborhood contains all trees within distance $k$ for several classic distance functions. For the nearest neighbor interchange (NNI), subtree prune and regraft (SPR), tree bisection and reconnection (TBR), and edge contraction and refinement (ECR) distances, we show that, under the exponential time hypothesis, there are no algorithms with running time $|I| o(k)$ where $|I|$ is the total input size. Hence, brute-force algorithms with running time $|X|^{O(k)} \cdot |I|$ are essentially optimal.

In contrast to the above distances, we observe that for the sECR-distance, where the contracted edges are constrained to form a subtree, a better solution within distance $k$ can be found in $k^{O(1)}$ time.

2012 ACM Subject Classification Theory of computation → Parameterized complexity and exact algorithms; Applied computing → Molecular evolution; Applied computing → Computational genomics

Keywords and phrases phylogenetic trees, parameterized complexity, tree distances, NNI, TBR

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1 Introduction

Maximum Parsimony is one of the most popular methods for inferring phylogenetic (evolutionary) trees from sequences of morphological or molecular characters. Given sequences of characters for $n$ taxa, this method reconstructs a phylogenetic tree $T$ whose $n$ leaves are labeled bijectively by the $n$ taxa and that has the minimum parsimony score over all such trees. The parsimony score is the number of character state changes along the tree edges that are necessary when extending the sequences for the leaves of $T$ to all internal vertices of $T$. Note that for each character, this score is at least $s - 1$, where $s$ denotes the number
of different character states. A phylogenetic tree is called perfect if it achieves score \( s - 1 \) for every character. Such a perfect phylogeny does not always exists. For a more comprehensive introduction to Maximum Parsimony, we refer the interested reader to [9].

From an algorithmic point of view, the Maximum Parsimony problem is notoriously hard: It is NP-complete even for binary characters [12]. Moreover, the current best running time is \( \Omega((2n - 3)!!) \), where \( (2n - 3)!! = 1 \cdot 3 \cdot \ldots \cdot (2n - 5) \cdot (2n - 3) \) [4]. The associated algorithm generates all possible binary phylogenetic trees on \( n \) leaves in a bottom-up fashion. Hence, the best known algorithm is essentially a brute-force-method. This running time bound is impractical when \( n > 15 \). Better running times are possible when the instance has a near-perfect phylogeny and the number of different character states \( s \) is small. Here, the running time is measured also in terms of the excess \( q \) over the score of a perfect phylogeny. In the general case, Maximum Parsimony can be solved in \( nm^{O(q)}2^{O(q^2 s^2)} \) time [10], where \( m \) is the length of the character sequences. In 2007, the running time was improved to \( O(21^q + 8^q nm^2) \) for the special case of binary characters and the practical usefulness of the improved algorithm was demonstrated for \( q \leq 10 \) [30]. In the worst case, however, \( q \) can be essentially as large as \( m \). Moreover, Maximum Parsimony is NP-hard even for \( q = 0 \) when the number of different character states is unbounded [3].

Given the hardness of Maximum Parsimony, solving this problem exactly is impractical for many real-world datasets due to prohibitive running times. Consequently, heuristic approaches, in particular local search, play an important role in computing good, but not necessarily optimal, solutions [2, 13, 14, 16, 17, 18, 19, 25, 26]. These approaches search the space of all possible phylogenetic trees on \( n \) taxa. In the course of such a search, the parsimony score of a subset of the phylogenetic trees in the space is computed. For any given tree, this step takes polynomial time using Fitch’s or Sankoff’s algorithm [11, 28]. A search through tree space starts by first computing a starting tree \( T \) before computing the parsimony score of all neighbors of \( T \). If there is a neighboring tree \( T' \) whose parsimony score is smaller than that of \( T \), then the search is continued by computing the parsimony score of all neighbors of \( T' \) and so on until a local optimum is found. In each iteration of the search, the neighboring trees are those that can be obtained from the current best tree by one or more rearrangement operations. The most well-known rearrangement operations on trees that are also considered in local search approaches for Maximum Parsimony, are nearest neighbor interchange (NNI), subtree prune and regraft (SPR), and tree bisection and reconnection (TBR) [1]. Each of these operations deletes an edge of a tree and then reconnects the resulting two subtrees. Depending on the operation, the reconnection is more or less restrictive, with SPR being a generalization of NNI and TBR being a generalization of SPR. The set of all trees that can be obtained by one operation is called the NNI, SPR, or TBR neighborhood, respectively. More general, we say that a tree \( T' \) is in the \( k \)-neighborhood with respect to NNI, SPR, or TBR of another tree \( T \), if \( T' \) can be obtained from \( T \) by at most \( k \) NNI, SPR, or TBR operations, respectively.

In addition to NNI, SPR, and TBR, the \( k \)-ECR operation has also been considered in the literature (see for example the works by Ganapathy et al. [13, 14]). This latter operation first contracts up to \( k \) edges and then refines the resulting tree arbitrarily. Here, the \( k \)-ECR neighborhood contains all trees that can be obtained from a starting tree by applying one \( k \)-ECR operation. The 1-ECR neighborhood is exactly the NNI neighborhood, but the 2-ECR neighborhood strictly contains the set of trees reachable by two NNI moves [14]. The \( k \)-ECR neighborhood appeared earlier implicitly under the term sectorial search [19]. The \( k \)-sECR neighborhood, a restricted version of the \( k \)-ECR neighborhood where the contracted edges must form a subtree was considered by Sankoff et al. [29]. They found that for larger
values of $k$, the $k$-sECR neighborhood gives better results than the 1-ECR neighborhood or, equivalently, the NNI neighborhood. Guo et al. [20] found that exploring the $k$-ECR neighborhood is too costly and thus proposed a restriction of this neighborhood which already leads to very good local optima. Their approach contracts $k$ edges and then refines the resulting tree by using neighbor joining, a fast distance-based method to reconstruct phylogenetic trees. To summarize, local search is an important paradigm for designing heuristics for Maximum Parsimony, and it has been noted that larger neighborhoods such as the $k$-ECR neighborhood give better results at the cost of higher running times. So far, there is however no study of how hard exploring larger neighborhoods actually is.

To analyze the computational complexity of exploring neighborhoods under NNI, SPR, TBR, $k$-ECR, and $k$-sECR, we use the framework of parameterized local search [8, 15, 23, 24]. Here, one studies local search problems with a neighborhood whose size can be adjusted by a parameter $k$. In the canonical parameterized local search problem, one is then given some solution for an optimization problem and the question is whether there is a better solution in the $k$-neighborhood. Local search for any of the aforementioned neighborhoods that are associated with distances between two trees fits exactly into this framework: we are given a phylogenetic tree and want to know whether there is one with a better parsimony score in the $k$-neighborhood. Typically, the $k$-neighborhood has a size of $O(|I|^f(k))$, where $|I|$ is the input size. In our case, the input size $|I|$ is in $O(n^2 \cdot m)$. Thus, using a brute-force algorithm, one can find a better solution in the neighborhood if it exists in $|I|^f(k)$ time. The algorithmic question is now whether this can be done much faster. In particular, a running time of $f(k) \cdot |I|^{O(1)}$ would be desirable since the explosion in the running time would then depend only on $k$ and not on $|I|$. Parameterized algorithmics provides toolkits to design such algorithms or to show that such algorithms are unlikely. The latter can be done by showing W[1]-hardness with respect to $k$ [6, 7] or by giving tight running time bounds based on the exponential time hypothesis (ETH) [21].

Our results are as follows. We show that even when all characters are binary, searching the $k$-ECR neighborhood is W[1]-hard with respect to $k$. The reduction that we use to establish this result also shows that, under the ETH, a running time of $|I|^{O(k)}$ is necessary. Moreover, the reduction implies hardness for searching the $k$-neighborhood with respect to NNI, SPR, and TBR. In a nutshell, our results show that one cannot gain a substantial speed-up over the brute-force algorithm when trying to search these large neighborhoods. We then establish that $n^{O(k)} \cdot m$ time is sufficient to search the $k$-neighborhoods with respect to any of NNI, SPR, TBR, and $k$-ECR, giving tight upper and lower bounds for the running time dependence on $k$. Finally, we observe that the $k$-sECR neighborhood of Sankoff [29] can be searched in $k^{O(k)} \cdot |I|^{O(1)}$ time, making it possible to consider much larger values of $k$ than for the other neighborhoods. Let us remark that, while we formally study the decision problem that asks for the existence of a better tree in the $k$-neighborhood, our hardness results and algorithms also apply to the problem of finding an optimal tree in the $k$-neighborhood.

Proofs of statements marked with (*) are deferred to a full version of the article.

## 2 Preliminaries

For details about relevant definitions of parameterized complexity such as fixed-parameter tractability, W[1]-hardness, parameterized reductions and ETH, refer to the standard monographs [6, 7].
Graph notation. For a graph $G = (V,E)$ and a vertex set $K \subseteq V$, let $E(K)$ denote the set of edges of $G$ where both endpoints are from $K$. The subdivision of an edge $e \in E$ in $G$ results in the graph $G'$ obtained by removing $e$ from $G$ and adding a new vertex which is adjacent to both endpoints of $e$. Let $v$ be a vertex of degree 2 in $G$. The suppression of $v$ in $G$ results in the graph $G'$ obtained by removing $v$ from $G$ and joining both neighbors of $v$ by an edge.

Phylogenetic trees. Throughout this paper, $X$ denotes a non-empty finite set of taxa.

An unrooted phylogenetic $X$-tree (for short, $X$-tree) $T$ is a tree with leaf-set $X$ and where no vertex has degree 2. If all non-leaf vertices of $T$ have degree three, then $T$ is called binary. Furthermore, if an edge $e$ is incident with a leaf of $T$, then $e$ is called a pendant edge and, otherwise, an internal edge. For two disjoint sets of taxa $A$ and $B$, we say that $A|B$ is a split of an $X$-tree $T$ if there is an edge $e$ in $T$ such that the deletion of $e$ results in two subtrees where one has leaf set $A$ and the other has leaf set $B$. The set of all splits of $T$ is denoted by $\Sigma(T)$. Furthermore, we say that an $X$-tree $T'$ is a refinement of $T$ if $\Sigma(T) \subseteq \Sigma(T')$. Additionally, if $T'$ is binary, then $T'$ is a binary refinement of $T$. We say that two $X$-trees $T$ and $T'$ are isomorphic if $\Sigma(T) = \Sigma(T')$. Equivalently, two $X$-trees $T$ and $T'$ are isomorphic if there is a bijection $\varphi$ between the vertices of $T$ and the vertices of $T'$ such that $\varphi(x) = x$ for all $x \in X$, and for all distinct vertices $u$ and $v$ of $T$, $\{u,v\}$ is an edge of $T$ if and only if $\{\varphi(u), \varphi(v)\}$ is an edge of $T'$.

Now, let $T$ be an $X$-tree and let $V'$ be a subset of the vertices of $T$. Then $T(V')$ denotes the minimal subtree of $T$ containing all vertices in $V'$. Let $A$ be a non-empty and proper subset of $X$ and let $T$ be a binary $X$-tree. If $A|(X \setminus A)$ is a split of $T$, then the subtree $T(A)$ is a pendant $A$-tree. Moreover, the pseudo-root of $T(A)$ is the unique vertex of degree 2 in $T(A)$ if $|A| > 1$ and the unique vertex of $T(A)$, otherwise.

Maximum parsimony. A character\(^1\) $c$ on $X$ is a function $c : X \to C$. If $|C| = 2$, then $c$ is called a binary character. Intuitively, $C$ can be thought of as the underlying alphabet and each element in the alphabet is a character state. Let $T$ be an $X$-tree with vertex set $V$, and let $c$ be a character on $X$ whose set of character states is $C$. An extension $c^*$ of $c$ to $V$ is a function $c^* : V \to C$ such that $c^*(x) = c(x)$ for each taxon $x \in X$. Let $c^*$ be an extension of $c$. A mutation edge of $c^*$ in $T$ is an edge $\{u,v\}$ in $T$ such that $c^*(u) \neq c^*(v)$ and we let $\text{score}_{c^*}(T)$ denote the number of mutation edges of $c^*$ in $T$. Then the parsimony score of $c$ on $T$, denoted by $\text{score}_c(T)$, is obtained by minimizing $\text{score}_{c^*}(T)$ over all possible extensions $c^*$ of $c$. An extension $c^*$ that minimizes $\text{score}_{c^*}(T)$ is called an optimal extension of $c$ in $T$. Moreover the maximum parsimony score of $c$, denoted by $\text{MP}(c)$, is the parsimony score of $c$ minimized over all binary $X$-trees.

Now let $S = \langle c_1, c_2, \ldots, c_n \rangle$ be a sequence of characters on $X$. Then the parsimony score of $S$ on an $X$-tree $T$ is defined as $\text{score}_S(T) = \sum_{i=1}^{n} \text{score}_{c_i}(T)$, and similarly, the maximum parsimony score of $S$, denoted by $\text{MP}(S)$, is the parsimony score of $S$ minimized over all binary $X$-trees.

We may abuse notation by writing $c \in S$ if the character $c$ is contained in the sequence $S$.

**SPR and TBR.** Let $T$ be a binary $X$-tree. Let $e = \{u,v\}$ be an edge of $T$, and let $T_1$ and $T_2$ be the two trees obtained from $T$ by deleting $e$ and suppressing $u$ if its degree is 2. Without loss of generality, we may assume that $T_2$ contains $v$. If $T_1$ contains at least one

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\(^1\) Characters as defined here are not elements of some alphabet but functions that assign an element of some alphabet to each taxon.
We next define a generalization of the SPR operation. Again, let $T$ be a binary refinement of the tree obtained from $X$. We say that $T'$ has been obtained from $T$ by a single subtree prune and regraft (SPR) operation. We next define a generalization of the SPR operation. Again, let $e$ be an edge of $T$, and let $T_1$ and $T_2$ be the two trees obtained from $T$ by deleting $e$ and suppressing any resulting degree-2 vertices. For each $i \in \{1,2\}$, if $T_i$ has at least one edge, subdivide an edge in $T_i$ with a new vertex $v_i$ and, otherwise, set $v_i$ to be the single vertex of $T_i$. Obtain a binary $X$-tree $T'$ by adding the new edge $\{v_1, v_2\}$. We say that $T'$ has been obtained from $T$ by a single tree bisection and reconnection (TBR) operation.

**NNI, k-ECR, and k-sECR.** Let $T$ be a binary $X$-tree. Let $e = \{u, v\}$ be an edge of $T$ and $e' = \{v, w\}$ be an internal edge of $T$ that is adjacent to $e$. Let $T'$ be a binary $X$-tree obtained from $T$ by deleting $e$, suppressing $v$, subdividing an edge that is incident with $w$ with a new vertex $v'$, and joining $u$ and $v'$ via a new edge. We say that $T'$ has been obtained from $T$ by a single nearest neighbor interchange (NNI) operation. Equivalently, if $T'$ is a binary refinement of the tree obtained from $T$ by contracting $e'$ and $T'$ is non-isomorphic to $T$, then $T'$ is obtained from $T$ by a single NNI operation.

Now let $T$ be a binary $X$-tree, and let $k$ be a positive integer. Let $T'$ be a binary refinement of a tree obtained from $T$ by contracting $k$ (distinct) internal edges $E'$. If $T'$ and $T$ are non-isomorphic, then we say that $T'$ is a single $k$-edge contract and refine ($k$-ECR) operation [13] apart from $T$ and that $E'$ is a contraction set for $T$ and $T'$. Note that an NNI operation is a 1-ECR operation and vice versa. We denote the restricted version of a $k$-ECR operation that requires the $k$ contracted edges to form a subtree of $T$ as $k$-sECR [29].

**Distance measures.** Let $T$ and $T'$ be binary $X$-trees. For each $\Theta \in \{\text{NNI, SPR, TBR}\}$, the distance $d_{\Theta}(T, T')$ is defined as the minimum number of $\Theta$ operations to transform $T$ into $T'$ [1]. The distance $d_{\text{ECR}}(T, T')$ is defined as the smallest number $k$ such that $T$ and $T'$ are one $k$-ECR operation apart. Analogously, the distance $d_{\text{sECR}}(T, T')$ is defined as the smallest number $k$ such that $T$ and $T'$ are one $k$-sECR operation apart.

**Considered problems.** In this work, we consider the parameterized complexity of the following problem for each distance measure $d \in \{d_{\text{NNI}}, d_{\text{SPR}}, d_{\text{TBR}}, d_{\text{ECR}}, d_{\text{sECR}}\}$.

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**d-LS Maximum Parsimony**

**Input:** A set of taxa $X$, a binary $X$-tree $T$, a sequence of characters $S$, and an integer $k$.

**Question:** Is there a binary $X$-tree $T'$ with $d(T, T') \leq k$ and $\text{score}_S(T') < \text{score}_S(T)$?

### 3 Properties of the Considered Distance Measures

In this section, we analyze the relation of the different distance measures.

- **Observation 3.1** ([1, 27]). The distance measures $d_{\text{NNI}}, d_{\text{SPR}},$ and $d_{\text{TBR}}$ are metrics.

- **Lemma 3.2** (*). The distance measure $d_{\text{ECR}}$ is a metric.

- **Observation 3.3.** Let $T$ and $T'$ be distinct binary $X$-trees and let $k > 0$ be an integer. If $d_{\text{ECR}}(T, T') = k$, then there is a binary $X$-tree $\tilde{T}$ with $d_{\text{ECR}}(\tilde{T}, T') > 0$ such that $d_{\text{ECR}}(T, T') = d_{\text{ECR}}(T, \tilde{T}) + d_{\text{ECR}}(\tilde{T}, T')$. 

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The idea behind Observation 3.3 is to consider the connected components of $T$ induced by the contraction set $S$ between $T$ and $T'$. If $S$ forms a subtree of $T$, then $S$ is connected and $d_{ECR}(T, T') = d_{ECR}(T, T')$. Hence, the statement holds for $T = T'$. Otherwise, let $S$ be an inclusion-maximal subset of $S$, such that $\tilde{S}$ forms a subtree of $T$. Since $\tilde{S}$ is inclusion-maximal, we can obtain $T'$ from $T$ in two steps: First, we can obtain an intermediate $X$-tree $\tilde{T}$ from $T$ by an sECR operation with contraction set $\tilde{S}$. Second, we can obtain $T'$ from $\tilde{T}$ by an ECR operation with contraction set $S \setminus \tilde{S}$.

**Lemma 3.4 (⋆).** Let $T$ and $T'$ be binary $X$-trees. Then, $d_{NNI}(T, T') \geq d_{ECR}(T, T')$.

**Lemma 3.5.** Let $T$ and $T'$ be binary $X$-trees. Then, $d_{ECR}(T, T') \geq d_{SPR}(T, T')$.

**Proof.** Let $k = d_{ECR}(T, T')$. Hence, there is a set $S$ of $k$ internal edges in $T$ such that $T'$ can be obtained by an sECR operation with contraction set $S$. Let $V'$ be the vertices of $T$ incident with some edge of $S$ and let $V^*$ be the neighbors of $V'$ in $T$ that are not incident with any edge of $S$. Recall that by definition of sECR operations, the edges of $S$ induce a subtree of $T$. Hence, $T(V^*)$ is a binary $V^*$-tree having the set $S$ as internal edges. For each vertex $v$ of $V^*$, let $T_v$ denote the pendant subtree of $T$ with pseudo-root $v$ obtained by removing the edge between $v$ and the unique neighbor of $v$ in $V'$. Since $T'$ can be obtained by an sECR operation with contraction set $S$, $T'$ contains a subtree $T'_v$ isomorphic to $T_v$ for each vertex $v$ of $V^*$. Hence, $d_{SPR}(T, T') = d_{SPR}(T, T')$, where $T_S$ is obtained from $T$ by replacing $T_v$ by the auxiliary taxa $v$ for each vertex $v$ of $V^*$ and where $T'_S$ is obtained from $T'$ by replacing $T'_v$ by the auxiliary taxa $v$ for each vertex $v$ of $V^*$ [1]. Note that $T_S = T(V^*)$.

Hence, it remains to show that $d_{SPR}(T, T') \leq k$. Since $T$ is binary and the edges of $S$ induce a subtree of $T$, $|V^*| = |S| + 3$. Moreover, since for each set of taxa $X'$ and each two binary $X'$-trees $T$ and $T'$, $d_{SPR}(T, T') \leq |X'| - 3$ [1], we conclude $d_{SPR}(T, T') \leq |V^*| - 3 = |S| = k$. Consequently, $d_{SPR}(T, T') \leq k = d_{ECR}(T, T')$.

**Lemma 3.6 (⋆).** Let $T$ and $T'$ be binary $X$-trees. Then, $d_{ECR}(T, T') \geq d_{SPR}(T, T')$.

## 4 Hardness of $d$-LS Maximum Parsimony

In this section, we establish our main theorem.

**Theorem 4.1.** For each distance measure $d \in \{d_{NNI}, d_{ECR}, d_{SPR}, d_{TBR}\}$ and even if each character is binary, d-LS Maximum Parsimony is NP-complete, W[1]-hard when parameterized by $k$, and cannot be solved in $f(k) \cdot |I|^{o(k)}$ time for any computable function $f$, unless the ETH fails.

We reduce from CLIQUE which is NP-hard [22], W[1]-hard when parameterized by $k$ [7], and cannot be solved in $f(k) \cdot |I|^{o(k)}$ time for any computable function $f$, unless the ETH fails [5, 6].

**CLIQUE**

**Input:** An undirected graph $G = (V, E)$ and an integer $k$.

**Question:** Is there a clique of size $k$ in $G$, that is, a set of vertices $K$ of size $k$, such that $|E(K)| = \binom{k}{2}$?

Let $I = (G = (V, E), k)$ be an instance of CLIQUE and let $d \in \{d_{NNI}, d_{ECR}, d_{SPR}, d_{TBR}\}$ be a distance measure. We describe how to construct an equivalent instance $I' = (X, T = (V', E'), S, k')$ of $d$-LS Maximum Parsimony in polynomial time where $k' := k$ if $d \in \{d_{SPR}, d_{TBR}\}$ and $k' := 2k$ if $d \in \{d_{NNI}, d_{ECR}\}$.
We initialize \( S \) with the empty sequence.

For a vertex \( v \in V \), the pendant \( X_v \)-tree \( T_v \). The bold edges are the only edges of \( T_v \) that are not in \( R \).

The subtree \( q \) contains the seven vertices of \( X \) that are the only edges of \( T_v \) for each vertex \( v \in V \).

Definition of the characters of \( S \). Next, we define the characters of \( S \) which are all binary characters whose character states are 0 and 1. We obtain \( S \) by concatenating two sequences of characters, \( S_G \) and \( S_R \), which we describe in the following.

First, we describe the characters of \( S_G \). An overview of the characters is given in Table 1. We initialize \( S_G \) as the empty sequence.

\[ S_G := \text{the empty sequence.} \]

The subtree \( T \) connecting the pendant trees \( T_v \) for each vertex \( v \in V \).

**Definition of \( X \) and \( T \).** We start with an empty taxa set \( X \) and add for each vertex \( v \in V \), a set \( X_v \) consisting of the eight taxa:

\[ \text{in}^0_v, \text{in}^1_v, \text{in}^2_v, \text{in}^3_v, \text{out}^0_v, \text{out}^1_v, \text{out}^2_v, \text{out}^3_v \]

to \( X \). Additionally, we add a taxon \( x^* \) to \( X \). This completes the definition of \( X \).

Next, we define the binary \( X \)-tree \( T = (V', E') \). Since \( X \) contains \( 8 \cdot |V| + 1 \) taxa and each internal vertex of \( T \) has three neighbors, \( T' \) has \( 16 \cdot |V| \) vertices and \( 2 \cdot |X| - 3 = 16 \cdot |V| - 1 \) edges. By definition, \( V' \) is a superset of \( X \). Additionally, for each vertex \( v \in V \), the set \( V' \) contains the seven vertices:

\[ \text{in}^0_v, \text{in}^1_v, \text{in}^2_v, \text{in}^3_v, \text{out}^0_v, \text{out}^1_v, \text{out}^2_v, \text{out}^3_v, \text{r}^\text{in}_v, \text{r}^\text{out}_v, \text{r}^\text{mid}_v \]

The subtree \( T_v := T(X_v) \) is depicted in Figure 1a.

Moreover, \( V' \) contains \( |V| - 1 \) additional vertices \( q_i \) with \( i \in [2, |V|] \). Fix some arbitrary ordering of the vertices of \( V \) and let \( V(i) \) denote the \( i \)th vertex of \( V \) according to that ordering. The vertex \( q_2 \) is adjacent to \( r^\text{out}_{V(1)}, r^\text{out}_{V(2)} \), and \( q_3 \). For each \( i \in [3, |V| - 1] \), the vertex \( q_i \) is adjacent to \( q_{i-1}, q_{i+1} \), and \( r^\text{out}_{V(i)} \). Finally, \( q_{|V| - 1} \) is adjacent to \( q_{|V| - 1}, r^\text{out}_{V(|V|)} \), and \( x^* \). See Figure 1b for an illustration. This completes the definition of \( T \).

**Intuition.** The idea of the reduction is as follows: Some of the characters that we define in the following will ensure that each binary \( X \)-tree \( T' \) that improves over \( T \) contains a pendant subtree \( T'(X_v) \) for each vertex \( v \in V \). Further characters will ensure that there are only two non-isomorphic trees for \( T'(X_v) \) which are depicted in Figure 1a and Figure 2. Intuitively, these two choices then function as a selection gadget for selecting vertex \( v \) as a vertex of the sought clique \( K \). The budget \( k' \) bounds how many such vertices can be selected. Finally, further characters will ensure that \( T' \) improves over \( T \) only if \( E(K) \) contains at least \( \binom{\frac{k}{2}}{2} \) edges.
For each edge $e \in E$, we add a character $c_e$ to $S_G$. Let $e$ be an edge of $E$. We set $c_e(x^*) := 1$.

If $v$ is a vertex of $V$. If $v$ is an endpoint of $e$, we set $c_e(x) := 1$ for each taxon $x \in \{v^0, v^1, v^0, v^1\}$ and we set $c_e(x) := 0$ for each taxon $x \in \{\{v^0, v^1, v^0, v^1\}\}$. Otherwise, if $v$ is not an endpoint of $e$, we set $c_e(x) := 1$ for each taxon $x \in \{v^0, v^1, v^0, v^1\}$ and we set $c_e(x) := 0$ for each taxon $x \in \{\{v^0, v^1, v^0, v^1\}\}$. Let $S_G$ denote the sequence of characters $c_e$ for each edge $e \in E$.

Next, we define a character $c_{\text{mal}}$. We set $c_{\text{mal}}(x^*) := 1$. For each vertex $v \in V$, we set $c_{\text{mal}}(v^0) = c_{\text{mal}}(v^1) := 1$ and we set $c_{\text{mal}}(x) := 0$ for each taxon $x \in X_v \setminus \{v^0, v^1\}$. We add a sequence $S_{\text{mal}}$ of $(k_i - 1)$ copies of $c_{\text{mal}}$ to $S_G$. Intuitively, in a binary $X$-tree $T'$, if both endpoints of an edge $e \in E$ are in the selected set $K$, then the parsimony score of $c_e$ in $T'$ is exactly the parsimony score of $c_e$ in $T$ minus one. Moreover, if $T'$ is non-isomorphic to $T$, then the parsimony score of $S_{\text{mal}}$ in $T'$ is exactly the parsimony score of $S_{\text{mal}}$ in $T$ plus $|S_{\text{mal}}|$.

Hence, the characters of $S_{\text{mal}}$ act as a hurdle to ensure that $E(K)$ contains at least $|S_{\text{mal}}| + 1 = \binom{k_i}{2}$ edges.

Finally, for each vertex $v \in V$, we define four characters $c_{\text{v, in}}, c_{\text{v, out}}, c_{\text{v, ri}},$ and $c_{\text{v, ro}}$. For each taxon $x$ of $X \setminus X_v$, we set $c_{\text{v, in}}(x) := c_{\text{v, out}}(x) := c_{\text{v, ri}}(x) := c_{\text{v, ro}}(x) := 1$. Now, let $x$ be a taxon of $X_v$.

- If $x$ is in $\{v^0, v^1\}$, we set $c_{\text{v, in}}(x) := 1, c_{\text{v, out}}(x) := 0, c_{\text{v, ri}}(x) := 1, c_{\text{v, ro}}(x) := 0$.
- If $x$ is in $\{v^0, v^1\}$, we set $c_{\text{v, in}}(x) := 1, c_{\text{v, out}}(x) := 0, c_{\text{v, ri}}(x) := 1, c_{\text{v, ro}}(x) := 0$.
- If $x$ is in $\{v^0, v^1\}$, we set $c_{\text{v, in}}(x) := 0, c_{\text{v, out}}(x) := 1, c_{\text{v, ri}}(x) := 0, c_{\text{v, ro}}(x) := 0$.
- If $x$ is in $\{v^0, v^1\}$, we set $c_{\text{v, in}}(x) := 0, c_{\text{v, out}}(x) := 1, c_{\text{v, ri}}(x) := 0, c_{\text{v, ro}}(x) := 1$.

Let $\alpha := 2|X| \cdot (|K| + \binom{k_i}{2})$. Note that $\alpha$ is larger than $\text{score}(T') + \text{score}_{\text{mal}}(T')$ of any binary $X$-tree $T'$, since such a tree $T'$ contains less than $2|X|$ edges and $|S_E| + |S_{\text{mal}}| = |E| + \binom{k_i}{2} - 1$. For each vertex $v \in V$, we extend $S_G$ by

- a sequence $S_{\text{v, in}}$ of $\alpha$ copies of $c_{\text{v, in}}$,
- a sequence $S_{\text{v, out}}$ of $\alpha$ copies of $c_{\text{v, out}}$,
- a sequence $S_{\text{v, ri}}$ of $2\alpha$ copies of $c_{\text{v, ri}}$, and
- a sequence $S_{\text{v, ro}}$ of $2\alpha$ copies of $c_{\text{v, ro}}$.

Let $S_v$ denote the combined sequences of $S_{\text{v, in}}, S_{\text{v, out}}, S_{\text{v, ri}},$ and $S_{\text{v, ro}}$. Intuitively, for each binary $X$-tree $T'$ that improves over $T$ and contains $T'(X_v)$ as a pendant subtree, the characters of $S_v$ ensure that $T'(X_v)$ is isomorphic to either the pendant tree depicted in Figure 1a or the pendant tree depicted in Figure 2. These two choices then function as a selection gadget for the vertices of the sought clique in $G$. This completes the construction of $S_G$. Note that $|S_G| = |E| + \binom{k_i}{2} - 1 + 6\alpha \cdot |V|$.

Next, we describe the sequence of characters $S_R$. Let $\beta := 2|X| \cdot |S_G|$. Note that $\beta$ is larger than $\text{score}_{\text{R}}(T)$ of any binary $X$-tree $T$, since such a tree $T$ contains less than $2|X|$ edges. Let $R := E' \setminus \{(v, v), (v, v), (v, v) \mid v \in V\}$. For each edge $e$ of $R$, we define a character $c_{R}$. Let $A$ be the split of $T$ induced by $e$. For each taxon $x \in A$, we set $c_{R}(x) := 0$ and for each taxon $x \in B$, we set $c_{R}(x) := 1$. We add as sequence $S_{R}^0$ of $\beta$ copies of $c_{R}$ to $S_R$.

Intuitively, the characters of $S_R$ ensure that each binary $X$-tree $T'$ that improves over $T$, shares the split that is induced by $e$ in $T$ for each edge $e$ of $R$. This implies that $T'(X_v)$ is a pendant subtree of $T'$ for each vertex $v \in V$.

Properties of binary $X$-trees. Before we show the correctness of the reduction, we first make some observations about binary $X$-trees with the characters of the construction.

Note that for each binary $X$-tree $T'$ and each edge $e$ of $R$, $\text{score}_{\text{R}}(T') \geq 1$. 

Table 1 An overview of the characters of $S_G$.

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<tr>
<th>$c_v \in S_E$</th>
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<th>$c \in S_w$</th>
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<td>$v \in e$</td>
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Definition 4.2. Let $T'$ be a binary X-tree. We say that $T'$ is split-consistent for $T$ and $R$ if for each edge $e$ of $R$, the split of $T$ induced by $e$ is also a split of $T'$.

In preparation for the next observation, note that if a binary X-tree $T'$ is not split-consistent for $T$ and $R$, then there is some edge $e$ of $R$ such that $\text{score}_{S_E}(T') \geq 2$ and thus $\text{score}_{S_E}(T') \geq 2 \cdot \beta$. Hence, $\text{score}_{S_E}(T') \geq \text{score}_{S_R}(T') \geq \beta \cdot (|R| + \epsilon)$. Since $\beta > \text{score}_{S_E}(T)$, this implies $\text{score}_{S_E}(T') > \text{score}_{S_E}(T)$. Hence, we conclude the following.

Observation 4.3. Let $T'$ be a binary X-tree. a) If $\text{score}_{S}(T') \leq \text{score}_{S}(T)$, then $T'$ is split-consistent for $T$ and $R$. b) If $T'$ is split-consistent for $T$ and $R$, then $\text{score}_{S_R}(T') = \beta \cdot |R|$.

To determine whether $I'$ is a yes-instance of d-LS MAXIMUM PARSIMONY, we analyze the structure of binary X-trees $T'$ with $\text{score}_{S}(T') \leq \text{score}_{S}(T)$. Due to Observation 4.3, we only need to consider binary X-trees that are split-consistent for $T$ and $R$ in the following.

Let $v$ be a vertex of $V$ and let $T'$ be a binary X-tree which is split-consistent for $T$ and $R$. Since there is an edge $e_v$ in $T$ such that $e_v$ induces the split $X_v || (X \setminus X_v)$ in $T$ and $e_v$ is contained in $R$, $X_v || (X \setminus X_v)$ is a split in $T'$. Hence, $T'(X_v)$ is a pendant tree. Moreover, since all edges incident with $v_1$ are in $R$, we can assume that $v_1$ is the common neighbor of $e_{v, \text{in}}$ and $e_{v, \text{out}}$ in $T'$. Similarly, we may assume that $\overline{v}_1$ is the common neighbor of $\overline{e}_{v, \text{in}}$ and $\overline{e}_{v, \text{out}}$ in $T'$, $\overline{v}_1$ is the common neighbor of $\overline{e}_{v, \text{in}}$ and $\overline{e}_{v, \text{out}}$ in $T'$, and $\overline{v}_1$ is the common neighbor of $\overline{e}_{v, \text{in}}$ and $\overline{e}_{v, \text{out}}$ in $T'$.

Definition 4.4. Let $T'$ be a binary X-tree which is split-consistent for $T$ and $R$, let $v$ be a vertex of $V$, and let $r$ be the pseudo-root of the pendant tree $T'(X_v)$. We say that $T'(X_v)$ is an in-rooting of $T_v$ if $v_1$ is adjacent to $r$, $\overline{v}_1$ has distance $2$ to $r$, and both $\overline{v}_1$ and $\overline{v}_1$ have distance $3$ to $r$. Similarly, we say that $T'(X_v)$ is an out-rooting of $T_v$ if $v_1$ is adjacent to $r$, $\overline{v}_1$ has distance $2$ to $r$, and both $\overline{v}_1$ and $\overline{v}_1$ have distance $3$ to $r$.

Figure 1a shows an out-rooting of $T_v$ and Figure 2 shows an in-rooting of $T_v$.

Note that for each vertex $v$ of $V$, there is a unique in-rooting of $T_v$ with respect to isomorphism. Similarly, there is a unique out-rooting of $T_v$ with respect to isomorphism. Note that for each vertex $v \in V$, $T_v$ is an out-rooting of $T_v$. We call a binary X-tree $T'$ well-rooted if $T'$ is split-consistent for $T$ and $R$ and if for each vertex $v \in V$, $T'(X_v)$ is either an in-rooting or an out-rooting of $T_v$. Note that $T$ is well-rooted.

Lemma 4.5 (*). Let $T'$ be a binary X-tree which is split-consistent for $T$ and $R$ and let $v$ be a vertex of $V$. If $T'(X_v)$ is an in-rooting of $T_v$ or an out-rooting of $T_v$, then $\text{score}_{S_E}(T') = 9\alpha$. Otherwise, $\text{score}_{S_E}(T') \geq 10\alpha$.  

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Next, we describe for a given well-rooted binary X-tree $T'$ the maximum parsimony scores of $T'$ with respect to the characters of $S_E$ and $S_{\text{mal}}$. The idea is that in a well-rooted binary X-tree $T'$, for each edge $e = \{u, v\} \in E$ where $T'(X_u)$ is an in-rooting of $T_u$ and where $T'(X_v)$ is an in-rooting of $T_v$, the parsimony score of the character $c_e$ in $T'$ is exactly the parsimony score of the character $c_e$ in $T$ minus one. Moreover, if $T'(X_v)$ is an in-rooting of $T_v$ for at least one vertex $v \in V$, then the parsimony score of the characters of $S_{\text{mal}}$ in $T'$ is exactly the parsimony score of the characters of $S_{\text{mal}}$ in $T$ plus $\binom{k}{2} - 1$.

Lemma 4.6. Let $T'$ be a well-rooted binary X-tree. Let $e = \{u, v\}$ be an edge of $E$.

(a) If $T'(X_u)$ is an in-rooting of $T_u$ and $T'(X_v)$ is an in-rooting of $T_v$, then $\text{score}_{c_e}(T') = 4(|V| - 2) + 2$. Otherwise, $\text{score}_{c_e}(T') = 4(|V| - 2) + 3$.

(b) If there is a vertex $w \in V$ such that $T'(X_w)$ is an in-rooting of $T_w$, then $\text{score}_{c_{\text{mal}}}(T') = |V| + 1$. Otherwise, that is, if $T'$ is isomorphic to $T$, $\text{score}_{c_{\text{mal}}}(T') = |V|$.

Proof. For each vertex $w$ of $V$, let $T'_w := T'(X_w)$. Let $V_{\text{in}}$ be those vertices $w$ of $V$, where $T'_w$ is an in-rooting of $T_w$ and let $V_{\text{out}} = V \setminus V_{\text{in}}$ be those vertices $w$ of $V$, where $T'_w$ is an out-rooting of $T_w$. For each vertex $w \in V_{\text{in}}$, let $r^m_w$ and $r^{\text{out}}_w$ be the neighbors of $r_{\text{mid}}_w$, and let $r_{\text{mid}}^w$ and $r_{\text{out}}^w$ be the neighbors of $r_{\text{mid}}^w$. Analogously, for each vertex $w \in V_{\text{out}}$, let $r_{\text{mid}}^w$ and $r_{\text{out}}^w$ be the name of the pseudo-root of $T'_w$, let $r_{\text{mid}}^w$ and $r_{\text{out}}^w$ be the neighbors of $r_{\text{mid}}^w$, and let $r_{\text{out}}^w$ and $r_{\text{mid}}^w$ be the neighbors of $r_{\text{mid}}^w$. Recall that since $T'$ is well-rooted, for each vertex $w \in V$, $r_{\text{mid}}^w$ is adjacent to both $r_{\text{mid}}^w$ and $r_{\text{out}}^w$, $r_{\text{out}}^w$ is adjacent to both $r_{\text{out}}^w$ and $r_{\text{mid}}^w$, and $r_{\text{out}}^w$ is adjacent to both $r_{\text{out}}^w$ and $r_{\text{mid}}^w$.

First, we show statement a). Let $c_e$ be a character for some edge $e = \{u, v\}$ of $E$. For each vertex $w \in V \setminus \{u, v\}$,

- let $P_{\text{in}}^w$ be the unique path between $r_{\text{mid}}^w$ and $r_{\text{out}}^w$ in $T'$,
- let $P_{\text{mid}}^w$ be the unique path between $r_{\text{mid}}^w$ and $r_{\text{out}}^w$ in $T'$,
- let $P_{\text{out}}^w$ be the unique path between $r_{\text{mid}}^w$ and $r_{\text{out}}^w$ in $T'$, and
- let $P_{\text{out}}^w$ be the unique path between $r_{\text{mid}}^w$ and $r_{\text{out}}^w$ in $T'$.

Note that each of these four paths only contains two edges and that these four paths are pairwise edge-disjoint. Let $P_w := \{P_{\text{in}}^w, P_{\text{mid}}^w, P_{\text{out}}^w, P_{\text{out}}^w\}$. Let $P'$ be a path in $P_w$ and let $w^0$ and $w^1$ be the terminals of $P'$. Since by definition $c_e(w^0) \neq c_e(w^1)$, for each extension $c_{e'}$ of $c_e$ in $T'$ at least one edge of $P'$ is a mutation edge of $c_{e'}$. Note that each path in $P_w$ is edge-disjoint with each path in $P_{w'}$ for distinct vertices $w$ and $w'$ of $V \setminus \{u, v\}$. Moreover, let $P_u$ be the path between $r_{\text{mid}}^u$ and $r_{\text{out}}^u$ in $T'$ and let $P_v$ be the path between $r_{\text{mid}}^v$ and $r_{\text{out}}^v$ in $T'$. Note that $P_u$ and $P_v$ are edge-disjoint and that both are edge-disjoint with each path $P_w \in P_w$ for each vertex $w \in V \setminus \{u, v\}$. Since $c_e(r_{\text{mid}}^u) = 0$ and $c_e(r_{\text{out}}^u) = 1$, for
each extension \( c^*_e \) of \( c_e \) in \( T' \), at least one edge of \( P_u \) is a mutation edge of \( c^*_e \). Similarly, since \( c_e(\overline{\text{in}}) = 0 \) and \( c_e(\text{out}) = 1 \), for each extension \( c^*_e \) of \( c_e \) in \( T' \), at least one edge of \( P_v \) is a mutation edge of \( c^*_e \). Hence, \( \text{score}_{c_e}(T') \geq 4(|V| - 2) + 2 \).

**Case 1**: \( T'_u \) is an in-rooting of \( T_u \) and \( T'_v \) is an in-rooting of \( T_v \). We define an extension \( c^*_e \) of \( c_e \) in \( T' \), such that \( \text{score}_{c^*_e}(T') = 4(|V| - 2) + 2 \). We set \( c^*_e(\text{out}) := c^*_e(\overline{\text{out}}) := c^*_e(\text{in}) := 0 \) and \( c^*_e(\overline{\text{in}}) := c^*_e(\text{out}) := c^*_e(\text{in}) := 0 \). For each remaining internal vertex \( v' \) of \( T' \), we set \( c^*_e(v') := 1 \). Hence, the edge set

\[
\{\{\text{out}, r_u\mid \text{mid}\}, \{\text{in}, r_v\mid \text{mid}\}\}
\]

contains the mutation edges of \( c^*_e \) in \( T' \). Consequently, \( \text{score}_{c^*_e}(T') = 4(|V| - 2) + 2 \) which implies \( \text{score}_{c_e}(T') \geq 4(|V| - 2) + 2 \).

**Case 2**: \( T'_u \) is an out-rooting of \( T_u \) or \( T'_v \) is an out-rooting of \( T_v \). Assume without loss of generality that \( T'_u \) is an out-rooting of \( T_u \). Let \( P_x^* \) be the unique path between \( \overline{\text{out}}_v \) and \( x^* \) in \( T' \). Since \( c_e(\overline{\text{out}}_v) = 0 \) and \( c_e(x^*) = 1 \), for each extension \( c^*_e \) of \( c_e \) in \( T' \), at least one edge of \( P_x^* \) is a mutation edge of \( c^*_e \). Note that \( P_x^* \) is edge-disjoint with \( P_u \) and edge-disjoint with each path \( P_w \) for each vertex \( w \in V \setminus \{u, v\} \). Moreover, since \( T'_u \) is an out-rooting of \( T_u \), \( P_x^* \) is also edge-disjoint with \( P_v \). Hence, \( \text{score}_{c_e}(T') \geq 4(|V| - 2) + 3 \). We define an extension \( c^*_e \) of \( c_e \) in \( T' \), such that \( \text{score}_{c^*_e}(T') = 4(|V| - 2) + 3 \). To this end, we distinguish whether \( T'_u \) is an in-rooting of \( T_u \) or an out-rooting of \( T_u \).

**Case 2.1**: \( T'_u \) is an in-rooting of \( T_u \). We set \( c^*_e(\overline{\text{out}}) := c^*_e(\overline{\text{in}}) := c^*_e(\text{in}) := 0 \) and \( c^*_e(\overline{\text{in}}) := c^*_e(\overline{\text{out}}) := 0 \). For each remaining internal vertex \( v' \) of \( T' \), we set \( c^*_e(v') := 1 \). Hence, the edge set

\[
\{\{\overline{\text{out}}, r_u\mid \text{out}\}, \{\overline{\text{in}}, r_v\mid \text{in}\}\}
\]

contains the mutation edges of \( c^*_e \) in \( T' \).

**Case 2.2**: \( T'_u \) is an out-rooting of \( T_u \). We set \( c^*_e(\overline{\text{in}}) := c^*_e(\overline{\text{in}}) := c^*_e(\text{in}) := 1 \) and \( c^*_e(\overline{\text{in}}) := c^*_e(\overline{\text{out}}) := 1 \). For each remaining internal vertex \( v' \) of \( T' \), we set \( c^*_e(v') := 0 \). Hence, the edge set

\[
\{\{\overline{\text{in}}, r_u\mid \text{in}\}, \{\overline{\text{in}}, r_v\mid \text{in}\}, \{\text{in}, q_n\}\}
\]

contains the mutation edges of \( c^*_e \) in \( T' \).

Consequently, in both cases \( \text{score}_{c^*_e}(T') = 4(|V| - 2) + 3 \) which implies \( \text{score}_{c_e}(T') = 4(|V| - 2) + 3 \).

Next, we show statement b). Consider the character \( c_{\text{mal}} \). For each vertex \( v \in V \), let \( P_v \) be the unique path between \( \overline{\text{out}}_v \) and \( \text{out}_v \) in \( T' \). Since \( c_{\text{mal}}(\overline{\text{out}}_v) = 0 \) and \( c_{\text{mal}}(\text{out}_v) = 1 \), for each extension \( c^*_e \) of \( c_{\text{mal}} \) in \( T' \) at least one edge of \( P_v \) is a mutation edge of \( c^*_e \). Note that the paths \( P_v \) and \( P_w \) are edge-disjoint for distinct vertices \( v \) and \( w \) of \( V \). Hence, \( \text{score}_{c_{\text{mal}}}(T') \geq |V| \).
Case 1: There is some vertex \( v \in V \) such that \( T'_v \) is an in-rooting of \( T_v \). Let \( P'_v \) be
the unique path between \( in_v^0 \) and \( x^* \) in \( T' \). Since \( c_{mal}(in_v^0) = 0 \) and \( c_{mal}(x^*) = 1 \), for each
extension \( c_{mal}^* \) of \( c_{mal} \) in \( T' \), at least one edge of \( P'_v \) is a mutation edge of \( c_{mal}^* \). Note that \( P'_v \)
is edge-disjoint with \( P_v \) for each vertex \( w \in V \) distinct from \( v \). Moreover, since \( T'_v \) is an
in-rooting of \( T_v \), \( P'_v \) is also edge-disjoint with \( P_v \). Hence, \( score_{mal}(T'_v) \geq \vert V \vert + 1 \). We define
an extension \( c_{mal}^* ) of \( c_{mal} \) in \( T' \), such that \( score_{mal}(T') = \vert V \vert + 1 \). We set \( c_{mal}(out_v) := 1 \),
for each vertex \( w \in V \). For each remaining internal vertex \( v' \) of \( T' \), we set \( c_{mal}(v') := 0 \). Hence,
the edge set \( \{(w_v, x_v') \} \cup \{(v_v, v'_v) \mid v \in V \} \) contains the mutation edges of \( c_{mal}^* \)
in \( T' \). Consequently, \( score_{mal}(T') = \vert V \vert + 1 \) which implies \( score_{mal}(T') = \vert V \vert + 1 \).

Case 2: For each vertex \( v \in V \), \( T'_v \) is an out-rooting of \( T_v \). Hence, \( T' \) is isomorphic to \( T \).
We define an extension \( c_{mal}^* \) of \( c_{mal} \) in \( T' \), such that \( score_{mal}(T') = \vert V \vert \). We set \( c_{mal}(in_v) := \\
\vspace{2mm}
c_{mal}(out_v) := c_{mal}(r_v^-) := c_{mal}(r_v^+) := 0 \), for each vertex \( v \in V \). For each remaining internal vertex \( v' \) of \( T' \), we set \( c_{mal}(v') := 1 \). Hence, the edge set \( \{\{r_v^-, r_v^+ \} \mid v \in V \} \) contains
the mutation edges of \( c_{mal}^* \) in \( T' \). Consequently, \( score_{mal}(T') = \vert V \vert \) which implies
that \( score_{mal}(T') = \vert V \vert + 1 \).

The score of improving \( X \)-trees with respect to \( S \). Since \( T \) is well-rooted, and for each vertex \( v \in V \), \( T_v \) is an out-rooting of \( T_v \), Observation 4.3, Lemma 4.5, and Lemma 4.6 imply
the following.

\begin{itemize}
  \item \textbf{Corollary 4.7.} \( score_S(T) = |E| \cdot (4(\vert V \vert - 2) + 3) + (\left(\frac{k}{2}\right) - 1) \cdot \vert V \vert + \vert V \vert \cdot 9\alpha + \vert R \vert \cdot \beta \).
\end{itemize}

Note that by definition, \( \alpha = 2(8\vert V \vert + 1) \cdot ((\left(\frac{k}{2}\right) > |E| \cdot (4(\vert V \vert - 2) + 3) + (\left(\frac{k}{2}\right) - 1) \cdot \vert V \vert .
Hence, \( score_S(T) < \alpha \cdot (9\vert V \vert + 1) + \vert R \vert \cdot \beta \).

\begin{itemize}
  \item \textbf{Corollary 4.8.} Let \( T' \) be a binary \( X \)-tree with \( score_S(T') < score_S(T) \). Then, \( T' \) is
well-rooted.
\end{itemize}

\textbf{Proof.} Due to Observation 4.3, \( T' \) is split-consistent for \( T \) and \( R \) and \( score_{S_R}(T') = \vert R \vert \cdot \beta \).
Assume towards a contradiction that there is a vertex \( v \in V \) such that \( T'(x_v) \) is neither
an in-rooting of \( T_v \) nor an out-rooting of \( T_v \). Hence, Lemma 4.5 implies \( score_{S_R}(T') \geq 10\alpha \) and \( score_{S_R}(T') \geq 9\alpha \) for each vertex \( w \in V \ \setminus \{v\} \). Consequently, \( score_{S_R}(T') \geq 10\alpha + (\vert V \vert - 1) \cdot 9\alpha + \vert R \vert \cdot \beta = \alpha \cdot (9\vert V \vert + 1) + \vert R \vert \cdot \beta > score_S(T) \), a contradiction.

\textbf{Distances between well-rooted binary \( X \)-trees.} Next, we describe for each distance measure \( d \in \{d_{NNI}, d_{ECR}, d_{SPR}, d_{TBR}\} \) the distance between \( T \) and any other well-rooted binary \( X \)-tree \( T' \).

\begin{itemize}
  \item \textbf{Lemma 4.9.} Let \( T' \) be a binary and well-rooted \( X \)-tree. Moreover, let \( K \) be the set of
vertices of \( V \) such that \( T'(x_v) \) is an in-rooting of \( T_v \) for each vertex \( v \in V \) and \( T'(x_w) \) is
an out-rooting of \( T_w \) for each vertex \( w \in V \ \setminus \ K \). Then, \( d_{NNI}(T, T') = d_{ECR}(T, T') = 2 \cdot |K| \)
and \( d_{SPR}(T, T') = d_{TBR}(T, T') = |K| \).
\end{itemize}

\textbf{Proof.} First, we show that \( d_{NNI}(T, T') = d_{ECR}(T, T') = 2 \cdot |K| \). To this end, we show
that \( d_{NNI}(T, T') \leq 2 \cdot |K| \) and that \( d_{ECR}(T, T') \geq 2 \cdot |K| \). Since \( d_{NNI}(T, T') \geq d_{ECR}(T, T') \)
due to Lemma 3.4, this then implies \( d_{NNI}(T, T') = d_{ECR}(T, T') = 2 \cdot |K| \).

To show that \( d_{NNI}(T, T') \leq 2 \cdot |K| \), we prove the following: Let \( T \) be a well-rooted binary \( X \)-tree
and let \( v \) be a vertex such that \( T(x_v) \) is an out-rooting of \( T_v \). Then, \( d_{NNI}(T, T) \leq 2 \),
where \( T \) is a well-rooted binary \( X \)-tree with \( T(X \ \setminus \ X_v) = T(X \ \setminus \ X_v) \) and where \( T(X_v) \)
We show that $\hat{T}$ with $A$ of $\in$-rooting of $d$ and that $\hat{T}$ is an in-rooting of the vertices of $\hat{T}(X_v)$ in $\hat{T}$. We obtain the well-rooted binary $X$-tree $\hat{T}$ from $\hat{T}$ by
- firstly removing the edges \{\text{in}_v, \text{out}_v\} and \{\text{mid}_v, r_v\} and adding the edges \{\text{mid}_v, r_v\} and \{\text{in}_v, \text{out}_v\}
- secondly removing the edges \{\text{in}_v, r_v\} and \{\text{mid}_v, r_v\} and adding the edges \{\text{mid}_v, r_v\} and \{\text{in}_v, \text{out}_v\}.

Since this can be done by two consecutive NNI operations and $\hat{T}(X \setminus X_v) = \hat{T}(X \setminus X_v)$, we conclude $d_{\text{NNI}}(\hat{T}, \hat{T}) \leq 2$. Since $d_{\text{NNI}}$ is a metric one can then show via induction over any arbitrary ordering of the vertices of $K$, that $d_{\text{NNI}}(T, T') \leq 2 \cdot |K|$.

It remains to show that $d_{\text{ECR}}(T, T') \geq 2 \cdot |K|$. Let $\hat{E}$ be a subset of the internal edges of $T'$, such that $T'$ can be obtained from $T$ by an ECR operation with contraction set $\hat{E}$. We show that $|\hat{E}| \geq 2 \cdot |K|$. Let $v$ be a vertex of $K$. Recall that $T_v$ is an out-rooting of $T_v$ and that $T_v'$ is an in-rooting of $T_v$. Hence, the edge $\{v, \text{mid}_v\}$ induces the split $A[B]$ in $T$ with $A := \{\text{in}_v, \text{in}_v^0, \text{out}_v, \text{out}_v^0, \text{mid}_v\}$ and $B := X \setminus A$. Since $A[B]$ is not a split of $T'$, the edge $\{\text{in}_v, \text{out}_v\}$ is contained in $E$. Similar, since the edge $\{\text{mid}_v, r_v\}$ induces the split $A[B]$ in $T$ with $A := \{\text{in}_v, \text{in}_v^1, \text{in}_v^0, \text{mid}_v\}$ and $B := X \setminus A$. Since $A[B]$ is not a split of $T'$, the edge $\{\text{mid}_v, r_v\}$ is contained in $E$. Hence, for each vertex $v$ of $V$, $\hat{E}$ contains at least two edges of $T(X_v)$. Consequently, $|\hat{E}| \geq 2 \cdot |K|$ which implies $d_{\text{ECR}}(T, T') \geq 2 \cdot |K|$.

Second, we show that $d_{\text{SPR}}(T, T') = d_{\text{TBR}}(T, T') = |K|$. Similar to the first part of the proof, we show that $d_{\text{SPR}}(T, T') \leq |K|$ and that $d_{\text{TBR}}(T, T') \geq |K|$. Since $d_{\text{SPR}}(T, T') \geq d_{\text{TBR}}(T, T')$ this then implies $d_{\text{SPR}}(T, T') = d_{\text{TBR}}(T, T') = |K|$.

To show that $d_{\text{SPR}}(T, T') \leq |K|$, we prove the following: Let $\hat{T}$ be a well-rooted binary $X$-tree and let $v$ be a vertex such that $\hat{T}(X_v)$ is an out-rooting of $T_v$. Then, $d_{\text{SPR}}(\hat{T}, \hat{T}) \leq 1$, where $\hat{T}$ is a well-rooted binary $X$-tree with $\hat{T}(X \setminus X_v) = \hat{T}(X \setminus X_v)$ and where $\hat{T}(X_v)$ is an in-rooting of $T_v$.
To show this claim, we describe an SPR operation transforming \( \hat{T} \) into \( \hat{T} \). See Figure 4 for an illustration of this SPR operation. Let \( r_v^\text{out} \) be the name of the pseudo-root of the pendant tree \( \hat{T}(X_v) \) and let \( q \) be the name of the unique neighbor of \( r_v^\text{out} \) outside of \( \hat{T}(X_v) \) in \( \hat{T} \). Moreover, let \( r_v^\text{in} \) be the name of the common neighbor of \( r_v \) and \( r_w \) in \( \hat{T} \). We obtain the well-rooted binary X-tree \( \hat{T} \) from \( T \) by: removing the edge \( \{r_v^\text{out}, q\} \), suppressing the vertex \( r_v^\text{out} \), subdividing the edge \( \{r_v^\text{in}, r_v^\text{out}\} \) by a vertex \( q' \), and adding the edge \( \{q, q'\} \). Since this can be done by a single SPR operation and \( \hat{T}(X \setminus X_v) = \hat{T}(X \setminus X_v) \), we conclude \( d_{\text{SPR}}(\hat{T}, \hat{T}) \leq 1 \). Since \( d_{\text{SPR}} \) is a metric, one can then show via induction over any arbitrary ordering of the vertices of \( K \), that \( d_{\text{SPR}}(T, T') \leq |K| \).

It remains to show that \( d_{\text{TBR}}(T, T') \geq |K| \). This proof is deferred to a full version of the article.

\[\triangleright\]

**Correctness.** Finally, we are able to show that \( I \) is a yes-instance of CLIQUE if and only if \( I' \) is a yes-instance of \( d \)-LS MAXIMUM PARSIMONY with appropriate distance bounds.

\[\triangleright\text{Lemma 4.10.} \text{ The following statements are equivalent:}\]

1. There is a clique of size \( k \) in \( G \).
2. There is a binary X-tree \( T' \) with \( \text{score}_S(T') < \text{score}_S(T) \) and \( d_{\text{SPR}}(T, T') \leq k \).
3. There is a binary X-tree \( T' \) with \( \text{score}_S(T') < \text{score}_S(T) \) and \( d_{\text{TBR}}(T, T') \leq k \).
4. There is a binary X-tree \( T' \) with \( \text{score}_S(T') < \text{score}_S(T) \) and \( d_{\text{NNI}}(T, T') \leq 2k \).
5. There is a binary X-tree \( T' \) with \( \text{score}_S(T') < \text{score}_S(T) \) and \( d_{\text{ECR}}(T, T') \leq 2k \).

**Proof.** First, we show that Item 1 implies each of Item 2–5. Let \( K \subseteq V \) be a clique of size \( k \) in \( G \). Further, let \( T' \) be a well-rooted binary X-tree such that for each vertex \( v \in K \), \( T'(X_v) \) is an in-rooting of \( T_v \), and for each vertex \( v \in V \setminus K \), \( T'(X_v) \) is an out-rooting of \( T_v \). Due to Lemma 4.9, \( d_{\text{SPR}}(T, T') = d_{\text{TBR}}(T, T') = k \) and \( d_{\text{NNI}}(T, T') = d_{\text{ECR}}(T, T') = 2k \). It remains to show that \( \text{score}_S(T') < \text{score}_S(T) \). Since \( T' \) is well-rooted, due to Observation 4.3, \( \text{score}_{\text{NNI}}(T') = |R| \cdot \beta \) and due to Lemma 4.5, for each vertex \( v \in V \), \( \text{score}_{\text{NNI}}(T') = 9\alpha \). Moreover, since \( K \) is non-empty, we obtain by Lemma 4.6, that \( \text{score}_{\text{NNI}}(T') = (\binom{k}{2} - 1) \cdot (|V| + 1) \). Since \( K \) is a clique in \( G \), \( |E(K)| = \binom{k}{2} \). Finally, by Lemma 4.6, for each edge \( e \in E(K) \), \( \text{score}_{\text{NNI}}(T') = 4(|V| - 2) + 2 \), and for each edge \( e \in E \setminus E(K) \), \( \text{score}_{\text{NNI}}(T') = 4(|V| - 2) + 3 \). We conclude

\[
\text{score}_S(T') = |E| \cdot (4(|V| - 2) + 2) + \left( \binom{k}{2} - 1 \right) \cdot |V| \cdot 9\alpha + |R| \cdot \beta \\
= |E| \cdot (4(|V| - 2) + 3) + \left( \binom{k}{2} - 1 \right) \cdot |V| \cdot 9\alpha + |R| \cdot \beta - 1 = \text{score}_S(T) - 1,
\]

due to Corollary 4.7. Hence, \( T' \) is a binary X-tree with \( \text{score}_S(T') < \text{score}_S(T) \), \( d_{\text{SPR}}(T, T') = d_{\text{TBR}}(T, T') = k \), and \( d_{\text{NNI}}(T, T') = d_{\text{ECR}}(T, T') = 2k \).

Second, we show that each of Item 2–5 implies Item 1. Let \( T' \) be a binary X-tree with a) \( \text{score}_S(T') < \text{score}_S(T) \) and b) \( d_{\text{SPR}}(T, T') \leq k \), \( d_{\text{TBR}}(T, T') \leq k \), \( d_{\text{NNI}}(T, T') \leq 2k \), or \( d_{\text{ECR}}(T, T') \leq 2k \). Since \( \text{score}_S(T') < \text{score}_S(T) \), due to Corollary 4.8, \( T' \) is well-rooted, that is, for each vertex \( v \in V \), \( T'_v := T'(X_v) \) is either an in-rooting of \( T_v \) or an out-rooting of \( T_v \). Let \( K \subseteq V \) be the set of all vertices \( v \) of \( V \) where \( T'_v \) is an in-rooting of \( T_v \). We show that \( K \) is a clique of size \( k \) in \( G \). Since \( d_{\text{SPR}}(T, T') \leq k \), \( d_{\text{TBR}}(T, T') \leq k \), \( d_{\text{NNI}}(T, T') \leq 2k \), or \( d_{\text{ECR}}(T, T') \leq 2k \), Lemma 4.9 implies that \( K \) is not empty. Moreover, since \( \text{score}_S(T') < \text{score}_S(T) \), \( T' \) is not isomorphic to \( T \), which implies that \( K \) is nonempty. Hence due to Lemma 4.6, \( \text{score}_{\text{NNI}}(T') = (\binom{k}{2} - 1) \cdot (|V| + 1) \). Moreover, since \( T' \) is well-rooted, due to Observation 4.3, \( \text{score}_{\text{NNI}}(T') = |R| \cdot \beta \) and due to Lemma 4.5,
for each vertex \( v \in V \), \( \text{score}_S(T') = 9a \). Finally, by Lemma 4.6, for each edge \( e \in E \setminus E(K) \), \( \text{score}_e(T') = 4(|V| - 2) + 3 \), and for each edge \( e \in E(K) \), \( \text{score}_e(T') = 4(|V| - 2) + 2 \). Consequently, \( \text{score}_S(T) - \text{score}_S(T') = |E(K)| - \binom{|K|}{2} - 1 \).

Since \( \text{score}_S(T') < \text{score}_S(T) \), we have \( |E(K)| \geq \binom{|K|}{2} \). Hence, \( K \) is a size-\( k \) clique in \( G \).

Since \( k' = k \) if \( d \in \{d_{	ext{SPR}}, d_{	ext{TBR}}\} \) and \( k' = 2k \) if \( d \in \{d_{	ext{NNI}}, d_{	ext{ECR}}\} \), Lemma 4.10 implies that \( I \) is a yes-instance of \( \text{CLIQUE} \) if and only if \( I' \) is a yes-instance of \( d-\text{LS MAXIMUM PARSIMONY} \). This completes the proof of Theorem 4.1.

## 5 Essentially Tight Brute-Force Algorithms

We now show that simple brute-force algorithms for \( d-\text{LS MAXIMUM PARSIMONY} \) for each distance measure \( d \in \{d_{	ext{NNI}}, d_{	ext{ECR}}, d_{	ext{SPR}}, d_{	ext{TBR}}\} \) essentially match the lower bounds shown in Theorem 4.1. First, consider a distance measure \( d \in \{d_{	ext{NNI}}, d_{	ext{SPR}}, d_{	ext{TBR}}\} \).

### Observation 5.1

Let \( T \) be a binary X-tree, let \( d \in \{d_{	ext{NNI}}, d_{	ext{SPR}}, d_{	ext{TBR}}\} \) be a distance measure, and let \( k \) be an integer. One can enumerate all binary X-trees \( T' \) with \( d(T, T') \leq k \) in \( |X|^{\Theta(k)} \) time.

Observation 5.1 can be seen as follows: there are \( |X|^{\Theta(1)} \) many binary X-trees \( T' \) such that \( d(T, T') = 1 \), all these trees can be enumerated in \( |X|^{\Theta(1)} \) time, and for each binary X-tree \( T' \) with \( d(T, T') > 0 \), there is a binary X-tree \( \hat{T} \) with \( d(\hat{T}, T') = 1 \) and \( d(T, T') = d(T, \hat{T}) + 1 \).

Furthermore, we may enumerate all binary X-trees \( T' \) with \( d_{	ext{ECR}}(T, T') \leq k \) as follows:

First, we enumerate all subtrees of \( T \) with at most \( k \) edges. Second, for each connected subtree \( T_s \) of \( T \) with at most \( k \) edges, we enumerate all binary refinements of \( T_s \) after contracting all edges of \( T_s \). In Lemma 5.2, we show that the first step can be done in \( \mathcal{O}(4^k \cdot k^{-0.5} \cdot |X|) \) time. In Lemma 5.3, we show that both steps can be performed in \( \mathcal{O}((2k + 1)!! \cdot 4^k \cdot \sqrt{k} \cdot |X|^2) \) time where \( (2k + 1)!! := 1 \cdot 3 \cdot \ldots \cdot (2k + 1) \).

### Lemma 5.2 (*)

For every binary X-tree \( T \) and every integer \( k \), all connected subtrees of \( T \) with at most \( k \) edges can be enumerated in \( \mathcal{O}(4^k \cdot k^{-0.5} \cdot |X|) \) time.

### Lemma 5.3 (*)

For a given binary X-tree \( T \) and an integer \( k \), there are \( \mathcal{O}((2k + 1)!! \cdot 4^k \cdot 2^{-0.5} \cdot |X|) \) binary X-trees \( T' \) with \( d_{	ext{ECR}}(T, T') \leq k \). Moreover, all these binary X-trees can be enumerated in \( \mathcal{O}((2k + 1)!! \cdot 4^k \cdot k \sqrt{k} \cdot |X|^2) \) time.

Hence, we obtain the following due to the fact that the parsimony score of a given X-tree can be computed in \( \mathcal{O}(|X| \cdot |S|) \) time [11].

### Theorem 5.4

\( d_{	ext{ECR}} \)-LS MAXIMUM PARSIMONY can be solved in \( \mathcal{O}((2k + 1)!! \cdot 4^k \cdot k \sqrt{k} \cdot |X|^2 \cdot |S|) = 2^{\Theta(k \log k)} \cdot |X|^2 \cdot |S| \) time.

Finally, we describe how to enumerate all binary X-trees \( T' \) with \( d_{	ext{ECR}}(T, T') \leq k \).

### Lemma 5.5

Let \( T \) be a binary X-tree and let \( k \) be an integer. One can enumerate all binary X-trees \( T' \) with \( d_{	ext{ECR}}(T, T') \leq k \) in \( |X|^{\Theta(k)} \) time.

**Proof.** We show this statement by induction over \( k \).

**Base case.** Consider \( k = 0 \). Hence, \( T \) is the only binary X-tree \( T' \) with \( d_{	ext{ECR}}(T, T') = 0 \) and can be enumerated in \( |X|^{\Theta(1)} \) time.
**Inductive step.** For the inductive step, suppose that for each binary X-tree $\tilde{T}$ and for each $k' < k$, one can compute all binary X-trees $T'$ with $d_{EKR}(\tilde{T}, T') \leq k'$ in $|X|^{O(k')}$ time. Note that this implies that for each $k' < k$ there are $|X|^{O(k')}$ binary X-trees $T'$ with $d_{EKR}(\tilde{T}, T') = k'$. For each $i < k$, let $T_i$ be the collection of all binary X-trees $T$ with $d_{EKR}(T, \tilde{T}) = i$ and let $T_{<k}$ be the collection of all binary X-trees $\tilde{T}$ with $d_{EKR}(T, \tilde{T}) < k$, that is, $T_{<k} = \bigcup_{i=0}^{k-1} T_i$. Moreover, let $T_{EKR}$ be the collection of all binary X-trees $\tilde{T}$ with $d_{EKR}(T, \tilde{T}) = k$. Note that $T_{<k}$ can be computed in $|X|^{O(k)}$ time and due to Lemma 5.3, $T_{EKR}$ can be computed in $k^{O(k)} \cdot |X|^{O(1)}$ time. Let

$$T_k := T_{EKR} \cup \bigcup_{i=1}^{k-1} \bigcup_{\tilde{T} \in T_i} \{T' \mid d_{EKR}(\tilde{T}, T') \leq k - i\}.$$ 

Recall that by the induction hypothesis, for each $i < k$, $T_i$ has size $|X|^{O(i)}$ and for each binary X-tree $\tilde{T} \in T_i$ the collection $\{T' \mid d_{EKR}(\tilde{T}, T') \leq k - i\}$ can be computed in $|X|^{O(k-i)}$ time. Hence, $T_k$ can be computed in $|X|^{O(k)}$ time. We set $T := T_k \cup T_{<k}$ and show that $T$ contains exactly the binary X-trees $T'$ with $d_{EKR}(T, T') \leq k$.

Assume towards a contradiction that this is not the case.

**Case 1:** There is a binary X-tree $T'$ with $d_{EKR}(T, T') \leq k$ such that $T'$ is not in $T$. By definition, $T_{<k}$ contains all binary X-trees $\tilde{T}$ with $d_{EKR}(T, \tilde{T}) < k$. Consequently, $d_{EKR}(T, T') = k$. Hence, due to Observation 3.3, there is a binary X-tree $\tilde{T}$ with $d_{EKR}(\tilde{T}, T') > 0$ such that $d_{EKR}(T, T') = d_{EKR}(T, \tilde{T}) + d_{EKR}(\tilde{T}, T')$. Let $i := d_{EKR}(T, \tilde{T})$.

Note that $i \leq k - 1$. If $i = 0$, then $T$ is isomorphic to $\tilde{T}$ and thus $d_{EKR}(T, T') = d_{EKR}(\tilde{T}, T') = k$. Hence, $T'$ is contained in $T_{EKR}$, a contradiction. Otherwise, if $i > 0$, then $\tilde{T}$ is contained in $T_i$. Moreover, since $d_{EKR}(\tilde{T}, T') = d_{EKR}(T, T') - d_{EKR}(T, \tilde{T}) = k - i$ and $d_{EKR}(\tilde{T}, T') \geq d_{EKR}(T, T')$, we have $d_{EKR}(\tilde{T}, T') \leq k - i$ which implies that $T'$ is contained in $T_i$, a contradiction.

**Case 2:** There is a binary X-tree $T'$ with $d_{EKR}(T, T') > k$ such that $T'$ is contained in $T$. Hence, $T'$ is contained in $T_k \setminus T_{EKR}$. That is, there is some $i$ with $1 \leq i \leq k$ and a binary X-tree $\tilde{T}$ in $T_i$ such that $d_{EKR}(T, \tilde{T}) \leq k - i$. Since $d_{EKR}$ is a metric, due to the triangle inequality, $d_{EKR}(T, T') \leq d_{EKR}(T, \tilde{T}) + d_{EKR}(\tilde{T}, T') \leq k$, a contradiction.

Since $T$ can be computed in $|X|^{O(k)}$ time, the statement holds.

We conclude the following.

**Theorem 5.6 (**). For each distance measure $d \in \{d_{NNI}, d_{EKR}, d_{SPR}, d_{TBR}\}$, $d$-LS Maximum Parsimony can be solved in $|X|^{O(k)} \cdot |S|$ time.

### 6 Conclusion

A clear goal for future research would be to improve the running time of the algorithm for the $k$-sEKR neighborhood. This seems promising since the current bottleneck is the enumeration of the binary refinements of the tree obtained after contracting $k$ edges. However, an algorithm for $d_{EKR}$-LS Maximum Parsimony running in $2^{n(k\log k)} \cdot |I|^{O(1)}$ time would imply an algorithm for Maximum Parsimony running in $2^{n(|X|\log |X|)} \cdot |I|^{O(1)}$ time: when applying the $d_{EKR}$-LS Maximum Parsimony algorithm with $k := |X| - 3$, locally optimal solution are also globally optimal. Hence, a more immediate question is whether Maximum
PARSIMONY can be solved in $2^{o(|X| \log |X|)} \cdot |I|^{O(1)}$ time. A further goal would be to find other neighborhoods for which $d$-LS MAXIMUM PARSIMONY can be solved in time $f(k) \cdot |I|^{O(1)}$. Finally, it is open whether better running times are possible when searching the neighborhood not for a better tree but for a perfect phylogeny, that is, for a tree where for each character, the parsimony score is equal to the number of character states minus one.

References

18:18 Parameterized Local Search for Maximum Parsimony


Abstract

A factorization of a string $S$ is a partition of $w$ into substrings $u_1, \ldots, u_k$ such that $S = u_1u_2 \cdots u_k$. Such a partition is called equality-free if no two factors are equal: $u_i \neq u_j, \forall i, j$ with $i \neq j$. The maximum equality-free factorization problem is to find for a given string $S$, the largest integer $k$ for which $S$ admits an equality-free factorization with $k$ factors.

Equality-free factorizations have lately received attention because of their applications in DNA self-assembly. The best approximation algorithm known for the problem is the natural greedy algorithm, that chooses iteratively from left to right the shortest factor that does not appear before. This algorithm has a $\sqrt{n}$ approximation ratio (SOFSEM 2020) and it is an open problem whether there is a better solution.

Our main result is to show that the natural greedy algorithm is a $\Theta(n^{1/4})$ approximation algorithm for the maximum equality-free factorization problem. Thus, we disprove one of the conjectures of Mincu and Popa (SOFSEM 2020) according to which the greedy algorithm is a $\Theta(\sqrt{n})$ approximation.

The most challenging part of the proof is to show that the greedy algorithm is an $O(n^{1/4})$ approximation. We obtain this algorithm via prefix free factor families, i.e., a set of non-overlapping factors of the string which are pairwise non-prefixes of each other. In the paper we show the relation between prefix free factor families and the maximum equality-free factorization. Moreover, as a byproduct we present another approximation algorithm that achieves an approximation ratio of $O(n^{1/4})$ that we believe is of independent interest and may lead to improved algorithms. We then show that the natural greedy algorithm has an approximation ratio that is $\Omega(n^{1/4})$ via a clever analysis which shows that the greedy algorithm is $\Theta(n^{1/4})$ for the maximum equality-free factorization problem.

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Introduction

A factorization of a string $S$ is a partition of $S$ into substrings $u_1, u_2, \ldots, u_k$ such that $S = u_1u_2 \ldots u_k$. Factorizations are central objects of study in stringology, a famous example being the Lempel-Ziv algorithm [14]. String factorizations have many other applications as we show next. For instance, finding an occurrence of a string $v$ in a text $T$ can be formulated as $T$ admitting a factorization $T = uwv$. Then, a string $v$ is a prefix of another string $T$ if $T = vw$ and it is a suffix of $T$ if $T = vw$. Moreover, many string problems can be seen as string factorization problems [9] such as: Shortest Common Superstring, Longest Common Subsequence and Shortest Common Supersequence, to name a few. Another example of a string factorization problem is the Minimum Common String Partition [6, 7], a problem concerned with identifying factorizations for two strings such that the sequence of factors for one string is the permutation of the other’s.

In this paper we focus on the equality-free factorization, a special case of string factorization in which all factors are distinct. The equality-free factorization is a restricted variant of a more famous problem, termed generalized function matching which has a long history starting from 1979 (see, e.g., [12] and the references therein for more details). In the generalized function matching the input consists of a text $t$ over an alphabet $\Sigma_1$ and a pattern $p = p_1p_2 \ldots p_m$ over an alphabet $\Sigma_2$. The goal is to find an injective function from $f : \Sigma_2 \rightarrow \Sigma_1^+$ such that $t = f(p_1)f(p_2) \ldots f(p_m)$. Thus, the maximum equality free factorization problem is a particular case of the generalized function matching in which all the characters of the pattern $p$ are distinct. In turn, generalized function matching is a particular case of string equations, which is a notoriously difficult problem (see, e.g., the JACM paper [13]). In fact, even the version which restricts character-to-character function matching is extremely difficult, see [1], as opposed to the more restricted parameterized matching [2, 3, 10] which is simpler. Thus, maximum equality free factorization is part of family of fundamental problems in stringology.

The maximum equality free factorization problem is also motivated by applications in DNA synthesis [4]. More specifically, it is possible to produce short DNA fragments that will self-assemble into the wanted DNA structure. However, to obtain the desired structure, it is required that no two fragments are identical. Since the fragments must be short, one approach is to split the target DNA sequence into as many distinct pieces as possible.

Previous work

The equality-free factorization problem was first introduced by Condon, Maňuch and Thachuk [4] where it was presented as the string partitioning problem. The string partitioning problem asks for a factorization into distinct factors such that each factor is at most of a certain length. The problem was studied in a more general setting where the measure of collision between two factors is either equality or one is a prefix/suffix of the other. Condon et al. showed that these variants are \textsc{NP}-complete. More recently, Fernau, Manea, Mereș and Schmid [5] presented a similar problem that imposes a lower bound on the number of factors instead of an upper bound on factor length. Fernau et al. showed that this variant is also \textsc{NP}-complete. Afterwards, Schmid [9] studied the Fixed-Parameter Tractability of the two problems.

The decision version of the problem, that is, given a string $S$ and an integer $k$, decide if there exists an equality free factorization of $S$ with at least $k$ factors, is termed MAXEFF-s (this is the notation of Schmid [9] and we decide to use it for the sake of consistency). The optimization version, in which we are given $S$ and the goal is to find an equality free
factorization with as many factors as possible, is termed OPTEFF-s. The acronyms for the two problems were introduced in the previous papers [5, 9, 11] and we will use them in our paper for consistency (OPT stands for optimization).

Mincu and Popa [11] study OPTEFF-s and another variant named Maximum Gapped Equality Free Factorization (OPTGEFF-s). In the latter it is not required that all the characters of the input strings are part of the factorization. That is, the goal is to find an equality free factorization using a maximum number of factors of a substring of the input string. More formally, a gapped factorization of string $S$ over alphabet $\Sigma$ is a tuple $(u_1, u_2, \ldots, u_k)$ such that $S = \alpha_0 u_1 \alpha_1 u_2 \alpha_2 \cdots \alpha_{k-1} u_k \alpha_k$, where $u_i \in \Sigma^+$ are the factors and $\alpha_i \in \Sigma^*$ are the gaps. OPTGEFF-s asks, for a given string $S$, to find the largest integer $k$ such that $S$ admits a gapped equality-free factorization of size $k$. In [11] a $2$-approximation for the OPTGEFF-s and a $\sqrt{n}$-approximation (where $n$ is the size of the input string) for the OPTEFF-s were shown. Moreover, it was conjectured [11] that the greedy approximation ratio is $\Omega(\sqrt{n})$. Grüttemeier et al. [8] show a randomized algorithm for solving the MAXEFF-s with running time $2^k \cdot k^{O(1)} + O(n)$.

Our results

As mentioned, the best-known approximation algorithm, the greedy algorithm, for the OPTEFF-s has an approximation ratio of $\sqrt{n}$ and it was conjectured that the greedy algorithm has an approximation ratio of $\Theta(\sqrt{n})$. In this paper, we show a better approximation algorithm for OPTEFF-s with ratio $O(n^{1/4})$. We then use this algorithm to show that the greedy algorithm has the same approximation ratio of $O(n^{1/4})$. Hence, this disproves the conjecture from [11] saying that the approximation ratio of the greedy algorithm is $\Omega(\sqrt{n})$.

The challenge is to show that the greedy algorithm has an approximation ratio of $O(n^{1/4})$. To get our approximation ratio we start with an (approximate) prefix free solution for the version with gaps. Then, we use the prefix free property to map the factors of a solution returned by the greedy algorithm to the aforementioned prefix free solution. Moreover, besides the greedy algorithm, we introduce another approximation algorithm for OPTEFF-s with an approximation ratio of $O(n^{1/4})$, that uses some interesting techniques and is of independent interest. We claim that our techniques give some key insights and perhaps open the path for better approximation algorithms for the problem.

Finally, we use a clever analysis to also show that the greedy algorithm cannot have an approximation ratio better than $O(n^{1/4})$ and, hence, the approximation ratio of the greedy algorithm is $\Theta(n^{1/4})$ for the maximum equality-free factorization problem.

2 The prefix-free property

For the OPTGEFF-s problem (the version of factorization with gaps), a $2$-approximation algorithm via a reduction from a scheduling problem was shown in [11]. A natural direction for proving an approximation algorithm for OPTEFF-s is to transform a factorization with gaps, obtained from the approximation algorithm of OPTGEFF-s, into a solution without gaps. However, it is difficult to transform a given factorization with gaps into a factorization without gaps with roughly the same number of factors.

In this section we show that it is possible to transform a special case of a factorization with gaps into a factorization without gaps. We introduce the notion of a prefix-free gapped factorization, which has an important role in our algorithm and might be of independent interest.
Definition 1. Let $n,k \in \mathbb{N}$, let $S$ be a string of length $n$ and let $F_k = \{S_1,S_2,\ldots,S_k\}$ be a set of non-overlapping factors of $S$ (possibly with gaps). $F$ is a prefix-free gapped factorization of $S$ if for all $i \neq j$, $S_i$ is not a prefix of $S_j$.

Given a prefix-free gapped factorization $F$ such that $|F| = k$, we prove that there is a transformation of $F$ into a factorization without gaps with the same number of factors $k$, since each factor $S_i$ can be extended until the next factor $S_{i+1}$ without colliding with another factor $S_j$.

Lemma 2. Let $n,k \in \mathbb{N}$ and let $S$ be a string of length $n$ with a prefix-free gapped factorization $F$ with $|F| = k$. Then, there is an equality free factorization for $S$ without gaps with at least $k$ factors.

Proof. Denote $S = T_0S_1T_1S_2T_2\ldots T_{k-1}S_kT_k$. Denote with $R_i = S_iT_i$. Note that for each $i \neq j$, $R_i$ and $R_j$ are not prefixes of each other because their prefixes are $S_i$ and $S_j$, respectively, which are not prefixes of each other.

Now, consider $S = T_0R_1R_2\ldots R_k$. If, for all $i$, $T_0 \neq R_i$, then we have an equality free $(k+1)$-factorization. Otherwise, there exists an $i$ such that $T_0 = R_i$. We distinguish two cases.

In the first case, if $i < k$, then we set $Q_i = R_iR_{i+1}$. Thus, $Q_i$ and all other factors $R_j$ are still not prefixes of each other. $T_0$, which equals $R_i$, is not a prefix of any other $R_j$ (because it equaled $R_i$) and is shorter than $Q_1$. Hence $S$ has a $k$-factorization $S = R_iR_1R_2\ldots R_{i-1}Q_iR_{i+2}\ldots R_k$.

In the second case, if $i = k$, then we set $Q_k = R_{k-1}R_k$ and using a similar argument as above we obtain a $k$-factorization $S = R_kR_1R_2\ldots R_{k-2}Q_k$.

3 An $O(n^{1/4})$-approximation algorithm

In this section we show an algorithm that has an $O(n^{1/4})$-approximation to OPTEFF-s.

Our algorithm (see Algorithm 1) is composed of two algorithms: a greedy algorithm, called Greedy1 (see Algorithm 3), which always yields an $\sqrt{n}$-approximation, and a new algorithm (Algorithm 2) which is described next. Algorithm 1 simply selects the better of the two algorithms and returns it.

Algorithm 1 An $O(n^{1/4})$-approximation algorithm for OPTEFF-s.

Input: String $S$.
1 $F \leftarrow$ Algorithm 2($S$);
2 $G \leftarrow$ Algorithm 3($S$);
3 if $|G| > |F|$ then
4 $\quad$ return $G$
5 return $F$

The basic idea behind Algorithm 2 is to find, for every fixed integer $1 \leq i \leq 2\sqrt{n}$, a greedy equality free gapped factorization of the input string in which every factor has length exactly $i$. The algorithm chooses from these gapped fixed length factorizations, the factorization with the most factors. Then, due to Lemma 2, we append to each of these factors the following adjacent (possibly empty) gap and we obtain an equality free factorization. See Algorithm 2 for more details.
Algorithm 2 Fixed length greedy factorization.

**Input:** String $S$.

1. $F \leftarrow \emptyset$;
2. for $i \leftarrow 1$ to $2\sqrt{n}$ do
3. $j \leftarrow 1, G \leftarrow \emptyset$;
4. while $j \leq n - i$ do
5. if $S[j..j + i - 1] \notin G$ then
6. $G \leftarrow G \cup \{S[j..j + i - 1]\}$;
7. $j \leftarrow j + i - 1$;
8. if $|G| > |F|$ then
9. $F \leftarrow G$;
10. Extend each factor $w_i \in F$ until factor $w_{i+1}$ (for the last factor, extend it until the end of $S$);
11. return $F$

Lemma 3. Algorithm 2 yields an equality-free factorization without gaps.

**Proof.** First, in the for loop, at each step, Algorithm 2 adds to $G$ only distinct substrings of $S$. Then, notice that for every two factors $w_1, w_2 \in G$, it holds that $w_1$ is not a prefix of $w_2$, since both $w_1$ and $w_2$ have the same length. Therefore, $G$ is a prefix-free gapped factorization, and due to Lemma 2, the factors are extended as in line 11 to have an equality-free factorization without gaps.

**Analysis**

Here we prove that when the optimal solution, denoted OPT, has “many” factors, Algorithm 2 returns a good approximation of OPT.

Formally, let $F$ be the factorization returned by Algorithm 2. Let $\alpha$ be $n/|OPT|$. Notice that $|OPT| = n/\alpha$. We claim that $|F| = \Omega(n/\alpha^2)$.

We first give an overview of the proof. First, we prove in Lemma 5 that there are at least $n/2\alpha$ short factors (of length at most $2\alpha$) in OPT. Then we prove in Lemma 6 that there are at least $\Omega(n/\alpha^2)$ factors of the exact same length in OPT. Next, we prove in Lemma 9 that the factorization $F$ returned in Algorithm 2 is a 2-approximation of optimal fixed length factorization (see Definition 7). Finally, we prove in Lemma 10 that $|F| = \Omega(n/\alpha^2)$.

Lemma 4. An $x$-short factor of $S$ is a factor of length $\leq x$. An $x$-long factor of $S$ is a factor of length $> x$. When $x$ is clear we will simply call them short factors and long factors.

Lemma 5. There exist at least $n/2\alpha$ factors in OPT that are $2\alpha$-short.

**Proof.** Let $LF$ denote the set of $2\alpha$-long factors in OPT and $SF$ denote set of the $2\alpha$-short factors in OPT. We will use an argument on $n$, the length of $S$. Each long factor must be, by definition, of length $\geq 2\alpha + 1$. Hence, by length arguments, $|LF| \cdot (2\alpha + 1) + |SF| \cdot 1 \leq n$ and, hence, $|LF| \leq n/(2\alpha + 1) - |SF|/(2\alpha + 1)$. On the other hand, since $|OPT| = n/\alpha$, we have that $|SF| = n/\alpha - |LF|$. Putting these two equations together yields that $|SF| = n/\alpha - |LF| \geq n/\alpha - n/(2\alpha + 1) + |SF|/(2\alpha + 1)$ and hence, $|SF| - |SF|/(2\alpha + 1) \geq n/\alpha - n/(2\alpha + 1)$ which in turn yields $2\alpha |SF|/(2\alpha + 1) \geq (n\alpha + n)/\alpha(2\alpha + 1)$. Hence, $2\alpha^2 |SF| \geq n\alpha + n$ and $|SF| \geq n/2\alpha + n/2\alpha^2 \geq n/2\alpha$. ▶
Next, we show that among the short factors, $\Omega(1/\alpha)$ fraction of them actually have exactly the same length.

**Lemma 6.** There exists an integer $\ell \leq 2\alpha$ such that there are at least $n/4\alpha^2$ short factors in OPT of length exactly $\ell$.

**Proof.** By Lemma 5, there are at least $n/2\alpha$ short factors in OPT. The average number of factors of each short length, is at least $n/2\alpha = n/4\alpha^2$. By the pigeonhole principle, there exists an integer $\ell \leq 2\alpha$ such that there are at least $n/4\alpha^2$ short factors in OPT of length exactly $\ell$. $\blacksquare$

Next we prove that Algorithm 2 is a constant approximation algorithm to the problem of gapped factorization with fixed lengths.

**Definition 7.** The Fixed-Length Maximum Gapped Equality-Free Factorization Size ($\text{FLOptGEFF-s}$) problem is defined as follows. For a given string $S$ and an integer $r$, find the largest integer $m$, such that $S$ admits a gapped equality-free factorization of size $m$ where all factors are of length $r$.

In [11], the problem of $\text{OPTGEFF-s}$ is reduced to the Job Interval Selection Problem with $k$ intervals ($\text{JISPk}$, see Theorem 8), which has a 2-approximation algorithm.

**Theorem 8 (restated from [11]).** Given $n$ jobs containing $k$ time intervals each, find the maximum number of intervals that can be selected such that (i) no two intervals intersect and (ii) at most one time interval is selected per job.

Analogously to [11], $\text{FLOptGEFF-s}$ is reducible as well to $\text{JISPk}$, and here we briefly show the reduction.

**Lemma 9.** Algorithm 2 is a 2-approximation for $\text{FLOptGEFF-s}$.

**Sketch Proof.** We construct an instance of $\text{JISPk}$ with $O(n)$ jobs from a string $S$ with $n$ characters. For each distinct substring of $S$ with length $r$, we create a job. For each substring $s$ we add $[a,b)$ as a time interval of $s$ for all occurrences $s = S[a,b]$ in $S$.

Since $\text{JISPk}$ has a 2-approximation algorithm, we have that $\text{FLOptGEFF-s}$ has a 2-approximation algorithm as well. Moreover, the algorithm that approximates $\text{FLOptGEFF-s}(S,r)$ for some string $S$ and integer $r$ is in fact the inner loop of Algorithm 2, on the iteration where $i = r$. $\blacksquare$

We are ready to prove a lower bound on the number of factors returned by Algorithm 2.

**Lemma 10.** Let $F$ be the factorization returned by Algorithm 2. Then, $|F| = \Omega(n/\alpha^2)$.

**Proof.** By Lemma 6, there exists $\ell$ such that there are at least $n/4\alpha^2$ short factors in OPT of length $\ell$.

Let $S_{\text{ALG}}^\ell$ be the number of factors of length $\ell$ produced by Algorithm 2, let $S_{\text{GEFF}}^\ell$ be $\text{FLOptGEFF-s}(S,\ell)$, and let $S_{\text{OPT}}^\ell$ be the number of factors of length $\ell$ in OPT.

By Lemma 9, there is a polynomial algorithm that is a 2-approximation of the number of occurrences of a factor of length $\ell$ in OPT. Moreover, the algorithm behind Lemma 9 is in fact the inner loop of Algorithm 2. Notice that $\ell \leq 2\alpha \leq 2n/|OPT| \leq 2\sqrt{n}$, since $|OPT| \geq \sqrt{n}$ and therefore there is an iteration where $i = \ell$. Then,

$$S_{\text{OPT}}^\ell/2 \leq S_{\text{GEFF}}^\ell/2 \leq S_{\text{ALG}}^\ell$$

where the first inequality is due to the definition of $\text{FLOptGEFF-s}$ and the second inequality is due to Lemma 9.
Hence, combining Lemma 6 and Lemma 9, for the iteration where \( i = \ell \) on line 9, \( |G| = S^\alpha_{\text{ALG}} \geq (n/4\alpha^2)/2 \). Since the number of factors returned by the algorithm is at least \( |G| \) (i.e. \( |F| \geq S^\alpha_{\text{ALG}} \)), we have that \( |F| = \Omega(n/\alpha^2) \).

As stated before, Algorithm 1 is composed by two algorithms, Algorithm 2 and Algorithm 3. Algorithm 3 was introduced in [11] as Greedy1 algorithm. In a nutshell, consider starting “at the left” of the string and adding the next shortest substring (distinct from the already selected factors) to the incumbent factorization at each step of the algorithm. See [11] for details.

\section{Algorithm 3 Greedy1.}

\textbf{Input:} String S.
1. \( j \leftarrow 1, F \leftarrow \emptyset; \)
2. for \( i \leftarrow 1 \) to \( n \) do
3. \hspace{1em} if \( S[j..i] \notin F \) then
4. \hspace{2em} \( F \leftarrow F \cup S[j..i]; \)
5. \hspace{1em} \( j \leftarrow i + 1; \)
6. \hspace{1em} Extend the last factor of \( F \) until the end of \( S; \)
7. return \( F \)

It was proven in [11] that Greedy1 yields an equality-free factorization. Moreover, they prove that Greedy1 produces at least \( \Omega(\sqrt{n}) \) factors.

\begin{theorem}
Algorithm 1 is an \( O(n^{1/4}) \)-approximation polynomial time algorithm for the OPTEFF-s problem.
\end{theorem}

\textbf{Proof.} Combining Greedy1 with Algorithm 2, we have an algorithm that produces at least \( \Omega(\max((n/\alpha^2), \sqrt{n})) \) factors. This gives an approximation ratio of \( O(\min(n/\alpha^2, (n/\alpha)/\sqrt{n})) = O(\min(n, \sqrt{n}/\alpha)) \), which is maximized at \( \alpha = \sqrt{n} \), i.e. \( \alpha = n^{1/4} \).

Finally, notice that the both Greedy1 and Algorithm 2 run in polynomial time of at most \( O(n^{1/3} \log n) \).

\section{The natural greedy algorithm is a \( \Theta(n^{1/4}) \)-approximation}

In this section we prove that Greedy1 itself achieves an approximation ratio of \( O(n^{1/4}) \).

\begin{lemma}
Greedy1 is a \( 2 \)-approximation of Algorithm 2.
\end{lemma}

\textbf{Proof.} Let \( S \) be a string, and let \( \ell \) be a positive integer. Let \( F_\ell \) be a fixed \( \ell \) length gapped factorization on \( S \) such that \( |F_\ell| = \text{FLOptGEFF-s}(S, \ell) \). Let \( F_G \) be the factorization that is the output of the Greedy1 algorithm. We show that \( |F_G| \geq |F_\ell|/2 \).

We map each factor of \( F_\ell \) to a factor of \( F_G \) as follows. Let \( f \in F_\ell \) be a factor in \( F_\ell \) and let \( i \leq j \) be two indices such that \( f = S[i..j] \). In \( F_G \), denote the factors that cover \( S[i] \) and \( S[j] \) as \( g_i \) and \( g_j \), respectively. If \( g_i \neq g_j \), map \( f \) to \( g_j \). If \( f \) is a suffix of \( g_j \), then also map \( f \) to \( g_j \). Otherwise, \( f \) is fully contained in a factor of \( F_G \) and \( f \) is not a suffix of \( g_j \). Then, it must be the case that there is a factor \( g_s \) in \( F_G \) such that the suffix of \( g_s \) is exactly \( f \), as otherwise Greedy1 would have cut the factor \( g_j \) right after index \( j \), but \( f \) is not a suffix of \( g_j \). Therefore, map \( f \) to \( g_s \) (if there are more than one factors with \( f \) as a suffix in \( F_G \), map to one of them arbitrarily).
Now, let $g$ be a factor in $F_G$, and let $\hat{i} \leq \hat{j}$ be two indices such that $g = S[\hat{i}, \hat{j}]$. We claim that there are at most two factors in $F_I$ that are mapped to $g$. There is at most one factor in $F_I$ that overlaps $S[\hat{i}]$, and therefore mapped to $g$ because of overlapping two factors in $F_G$. Moreover, since all the factors in $F_I$ have the same size, there is at most one factor in $F_I$ such that the suffix of $g$ is equal to the factor. Therefore, there are at most 2 factors in $F_I$ that are mapped to $g$. To conclude, there are at most $2 \cdot |F_G|$ factors in $F_I$, for every $\ell$.

Let $A$ be the factorization returned by Algorithm 2. There is an $\ell$ such that $|F_G| \geq |F_A|$. Since we proved that $|F_G| \geq |F_1|/2$ for every $\ell$, we also have that $|F_G| \geq |F_A|/2$.

Combining the above lemma with the Theorem 11, we conclude the following theorem.

> **Theorem 13.** The approximation ratio of Greedy1 is $O(n^{1/4})$.

**Proof.** By Lemma 12, Greedy1 is a constant approximation of Algorithm 2 and therefore Greedy1 is also a constant approximation of Algorithm 1 (that simply uses Algorithm 2 and Greedy1 and returns the maximum between them). Since by Theorem 11 Algorithm 1 is an $O(n^{1/4})$-approximation for OPTEFF-s we have that Greedy1 is also an $O(n^{1/4})$-approximation.

## 5 Tightness of Algorithm 1

In this section we prove that our analysis of Algorithm 1 is actually tight. We show that there is a case where both Greedy1 and Algorithm 2 have an approximation ratio that is at least $\Omega(n^{1/4})$.

Similar to [11], we define a string $S$ as follows. Let $n$ be a square of an even number, i.e. there is an integer $k$ such that $n = (2k)^2$. Let $\Sigma = \{x_1, x_2, \ldots, x_{\sqrt{n}}\}$ be an alphabet. We define variables $X_1, X_2, \ldots, X_{\sqrt{n}}$ such that for each variable $X_i$, define $X_i = x_1 x_2 \ldots x_i$. The string $S$ is defined as $S = X_1 X_2 \ldots X_{\sqrt{n}}$. Note that $|S| = \Theta(n)$.

> **Lemma 14.** There exists a factorization of $S$ with $\Omega(n^{3/4})$ factors.

**Proof.** We first factorize $S$ into $\Omega(n^{3/4})$ factors with gaps, and afterwards we get rid of the gaps. We factorize the variables $X_1 \ldots X_{\sqrt{n}/2 - 1}$ using only one factor. Then, the variable $X_{\sqrt{n}/2}$ is factorized into $x_1 x_2 \ldots x_{\sqrt{n}/2}$, for a total of $\sqrt{n}/2$ factors. At least $\sqrt{n}/2 - 1$ factors are produced by the variables $X_{\sqrt{n}/2+1} X_{\sqrt{n}/2+2}$ as follows. The variable $X_{\sqrt{n}/2+1}$ is factorized into $x_1 x_2 x_3 x_4 \ldots$, for a total of $\lceil |X_{\sqrt{n}/2+1}]/2 \rceil$ factors. The variable $X_{\sqrt{n}/2+2}$ is factorized into $x_2 x_3 x_4 x_5 \ldots$, also for a total of at least $\lceil |X_{\sqrt{n}/2+2}]/2 \rceil$ factors.

In general, at each iteration $i$, the algorithm produces factors of length $i$ using $i$ variables and $i$ offsets. Each variable is of length at least $\sqrt{n}/2$, therefore at least $i \cdot \lceil (\sqrt{n}/2)/i \rceil \geq \sqrt{n}/2 - i$ factors are produced by $i$ variables. For each iteration $i$, the $r$th variable $X_j$ of iteration $i$ produces factors of length $i$ starting at index $r$ with respect to the beginning of $X_j$. This procedure produces an equality free factorization with gaps.

There is a constant $c > 0$ such that there are $c \sqrt{n}/2 = cn^{1/4}$ iterations to the process. Therefore, at least

$$\sum_{i=1}^{cn^{1/4}} \sqrt{n}/2 - i \geq cn^{3/4}/2 - c^2 \sqrt{n} = \Omega(n^{3/4})$$

factors are produced in this process.
We are left with handling the gaps. Notice that there are two reasons for a gap to occur. First, (1) on iteration \(i\) and variable \(X_j\), we produce \(\lfloor |X_j|/i \rfloor\) factors, and \(|X_j| - i \lfloor |X_j|/i \rfloor > 0\), so we have a gap at the end of \(X_j\). Second, (2) on iteration \(i\) and the \(r\)th variable of the iteration \(X_j\), when \(X_j\) is not the first variable of iteration \(i\) \((r \neq 1)\), the factorization of \(X_j\) does not start from \(x_1\) but from \(x_r\).

For gaps of type 1, let \(X_j\) be a variable that has a gap at the end of \(X_j\). Then, if \(j \neq \sqrt{n}\), we extend the first factor of \(X_j+1\) backwards to close the gap. This extended factor is unique since there is only one instance of \(x_j x_1\) in \(S\). If \(j = \sqrt{n}\) and we are in the last variable, we extend the last factor of \(X_j\). This extended factor is unique since there is only one instance of this length in \(S\).

For gaps of type 2, let \(X_j\) be a variable that has a gap at the beginning of \(X_j\). Then, we extend the last factor of \(X_j-1\) forward to close the gap. This extended factor is unique since there is only one instance of \(x_{j-1} x_1\) in \(S\).

For gaps with both types 1 and 2, we just handle them as gaps with type 1.

On string \(S\), Greedy1 produces \(\Theta(\sqrt{n})\) factors. Hence, and by Lemma 14, we have the following corollary.

\textbf{Corollary 15.} The approximation ratio of Greedy1 is \(\Omega(n^{1/4})\).

On the string \(S\) described above, Algorithm 2 produces \(O(\sqrt{n})\) factors. To see this, let \(l\) be some length that is being observed in line 2 of Algorithm 2. There are (at most) \(\sqrt{n}\) \(x_1\)’s in string \(S\). Therefore there are at most \(\sqrt{n}\) factors containing \(x_1\). On the other hand, every factor that does not contain \(x_1\), must start in a unique character (since before the extension, every factor is of length exactly \(l\)). There are (at most) \(\sqrt{n}\) unique characters in \(S\). Therefore, there are at most \(\sqrt{n}\) factors not containing \(x_1\).

Hence, and by Lemma 14, we have the following corollary.

\textbf{Corollary 16.} The approximation ratio of Algorithm 2 is \(\Omega(n^{1/4})\).

Finally, since both lower bounds were based on the same case of string \(S\), we have the following corollary.

\textbf{Corollary 17.} The approximation ratio of Algorithm 1 is \(\Omega(n^{1/4})\).

\section{Conclusions and future work}

In this paper we disproved one of the conjectures of Mincu and Popa [11] and show that the natural greedy algorithm for the OPTEFF-s problem has a \(\Theta(n^{1/4})\)-approximation factor. It is, of course, a natural open question to improve the approximation ratio for OPTEFF-s using a different algorithm than the greedy. We believe that the key in succeeding to obtain such an algorithm is finding a better approximation algorithm for the case when the number of factors in an optimal solution is relatively small. Thus, the ideas introduced in Section 3, where we present an alternative \(O(n^{1/4})\) approximation algorithm, represent a promising direction.

\begin{thebibliography}{9}
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String Factorization via Prefix Free Families


Improving the Sensitivity of MinHash Through Hash-Value Analysis

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Abstract
MinHash sketching is an important algorithm for efficient document retrieval and bioinformatics. We show that the value of the matching MinHash codes convey additional information about the Jaccard similarity of $S$ and $T$ over and above the fact that the MinHash codes agree. This observation holds the potential to increase the sensitivity of minhash-based retrieval systems. We analyze the expected Jaccard similarity of two sets as a function of observing a matching MinHash value $a$ under a reasonable prior distribution on intersection set sizes, and present a practical approach to using MinHash values to improve the sensitivity of traditional Jaccard similarity estimation, based on the Kolmogorov-Smirnov statistical test for sample distributions. Experiments over a wide range of hash function counts and set similarities show a small but consistent improvement over chance at predicting over/under-estimation, yielding an average accuracy of 61% over the range of experiments.

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1 Introduction
MinHash sketching is an important algorithm for efficient document retrieval. It reduces a set $S$ of size $n$ to a smaller representation of size $m \ll n$ by applying $m$ distinct hash functions $h_1,\ldots,h_m$ to each of the $n$ elements of $S$, and identifies the smallest hash code for each $h_i$. This vector of minimum hash codes serves a sketch for the larger set $S$. A classical result [2, 3] shows that the probability that smallest hash codes of two sets $S$ and $T$ are equal is identical to $J(S,T)$, the Jaccard similarity of $S$ and $T$. Thus the fraction of matching MinHash codes represents an unbiased estimator of $J(S,T)$.

Hash code values, by definition, are not supposed to mean anything. They represent mappings of an item $x$ to a pseudorandom integer $h(x)$ for purpose of fast identity matching and retrieval. The relative values of $h(x)$ and $h(y)$ for items $x$ and $y$ have no special properties beyond that of $h(x) = h(y)$ likely implies that $x = y$ for the conventional hash functions as employed in algorithms such as MinHash.

But in this paper, we report a curious observation associated with MinHash. Suppose the MinHash values for sets $S = \{s_1,\ldots,s_n\}$ and $T = \{t_1,\ldots,t_n\}$ equal both a particular value, namely:

$$a = \min_{j=1}^{n} h_i(s_j) = \min_{j=1}^{n} h_i(t_j)$$

We shall show that the value of this matching MinHash value $a$ conveys additional information about the Jaccard similarity of $S$ and $T$ over and above the fact that the MinHash values agree.

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This observation holds the potential to increase the sensitivity of minhash-based retrieval systems. Our main results in this paper are:

- We explain why observing a larger matching MinHash value increases the expectation of high similarity between \( S \) and \( T \). Specifically, the expected value of a common MinHash value \( a \) for two \( n \)-element sets with intersection size \( i \) is \( N/(2n - i + 1) \), presuming the underlying hash function selects an integer from \([0...N]\) uniformly at random.

- We analyze the expected Jaccard similarity of two sets as a function of observing a matching MinHash value \( a \) under a reasonable prior distribution on intersection set sizes, specifically the case where pairs of \( n \)-element sets have equal probability of intersection size \( i \) for \( 1 \leq i \leq n \). Experimental results confirm a modest increase in the sensitivity of our hash-code weighted variant of MinHash over the original, over a range of set similarities and number of hash codes.

- We present a practical approach to using MinHash values to improve the sensitivity of traditional Jaccard similarity estimation, based on the Kolmogorov-Smirnov statistical test for sample distributions. Our techniques provide a supplemental signal suggesting whether the fraction of matching MinHashes is more likely to over-estimate or underestimate the true Jaccard similarity between two sets. Experiments over a wide range of hash counts \((k)\) and set similarities show a small but consistent improvement over chance, specifically an average accuracy of 61% over the range of experiments.

We believe that this orthogonal view of measuring Jaccard similarity through the value of matching MinHash codes is novel, and will inspire further interest. Although the practical improvement we have demonstrated is not large, we believe that better interpretations of the underlying statistics may yield better results.

This paper is organized as follows. We begin with a survey of the literature of MinHash and related techniques in Section 2. We provide intuition as to why the value of the matching MinHash value matters through a thought experiment in Section 3. We present our analysis of the expected intersection size as a function of MinHash value for an appropriate prior distribution in Section 4, and ways to combine this information into an estimate of Jaccard similarity in Section 5. An alternate approach to interpret the values of MinHash codes using the Kolmogorov-Smirnoff statistical test is presented in Section 6. Finally, we conclude with some open problems raised by our work.

## 2 Prior Work

Broder [2, 3] developed MinHash as a solution to identifying similar documents (represented as sets of shingles or substrings) within a large text corpus while avoiding the quadratic-time costs of explicitly comparing each pair of documents. A function \( h(x) \) is applied to each set element, mapping each element \( x \) to a pseudo-random integer. Each set \( S \) is represented by the minimum hash value among all its elements.

The Jaccard similarity \( J(S, T) \) of two sets \( S \) and \( T \) is defined as

\[
J(S, T) = \frac{|S \cup T|}{|S \cap T|}.
\]

For identical sets, \( J(S, T) = 1 \) while for disjoint sets, \( J(S, T) = 0 \). Broder [2, 3] observed that the probability that two sets \( S \) and \( T \) generate the same MinHash value exactly equals the Jaccard similarity of the two sets, \( J(S, T) \). The fraction of matching minimum hashes over \( k \) independent hash functions provides an unbiased estimator of \( J(S, T) \), with the variance of this estimate equal to \( J(S, T)(1 - J(S, T))/k \) [4]. Surveys of MinHash sketching include Cohen [9].
MinHash is one of the most important algorithms for web search and duplicate detection [12, 15, 17], social networks [8], and machine learning [7, 19]. More recently, it has been successfully applied to bioinformatics for sketching large DNA sequence datasets, starting from the seminal Mash software [21] and followed by related tools [5, 28, 1, 20]. Such applications motivate our efforts in this paper to increase the sensitivity of minimum hashing-based similarity measures.

Locality Sensitive Hashing (LSH) techniques seek to map similar data objects to the same hash codes with a higher probability than the dissimilar ones by adopting a family of randomized hash functions. Indyk et al. [16, 14, 10] introduced the notion of locality-sensitive hashing in the context of nearest-neighbor search and string similarity. MinHash can be viewed an instance of locality sensitive hashing. An extended survey of locality-sensitive hashing can be found in [24].

SimHash [6] is a LSH-based method which provides an unbiased estimator of the similarity between two vectors. Specifically, the probability that two vectors $u$ and $v$ generate the same SimHash value equals the Cosine similarity of $u$ and $v$. Henzinger [15] performed a large-scale comparison of MinHash and SimHash on detecting similar web pages, finding that a hybrid of the two approaches yielded the best results. Srivastava and Li [22] present analysis and experiments to suggests that MinHash is more sensitive than SimHash in regions of high similarity.

Each subset element is granted equal weight in traditional MinHash schemes, but this can be generalized in weighted MinHash, perhaps to reflect the TD-IDF values of each word. Weighted MinHash algorithms are surveyed in [25].

Finally, we mention another related sketching scheme named HyperLogLog [13] primarily designed for the task of estimating the number of distinct items in a stream, but also capable of estimating the cardinality of the union of two sets and therefore their Jaccard similarity. Several works proposed unifying combinations of the two sketches [11, 27]. Note that HyperLogLog has some relationship with our work, as it employs the idea of estimating the cardinality of a random set from its minimum value. However, a direct application of this idea to MinHash has not been made, to our knowledge.

### 3 Thought Experiment: Why MinHash Values Matter

We present the following thought experiment to illustrate how the actual value of matching MinHash codes provides information about Jaccard similarity. For a set $S$, let $M_h(S)$ denote its MinHash value under a given hash function $h$, i.e. $M_h(S) = \min_{s \in S} \{ h(s) \}$. We use the notation $M(S)$ when $h$ is irrelevant (but assumed fixed across sets).

Now consider following two “extreme” situations involving pairs of sets $S$ and $T$, each with $n$ elements:

1. $S$ and $T$ are identical. Hence by definition, the minimum hash values must match, so $a_1 = M_h(S) = M_h(T)$.
2. $S$ and $T$ intersect in only one element $x$, which happens to be the minimum value of both under $h$, so $a_2 = h(x) = M_h(S) = M_h(T)$.

Now, given just the two unlabeled values for $a_1$ and $a_2$, can we correctly assign these codes to the appropriate case above with probability greater than $1/2$?

Assume hash function $h$ selects an integer from the range $[0 \ldots N]$ uniformly at random. Now suppose that two $n$-element sets with intersection size $i$ share a common MinHash value $m$. Then $m$ is the smallest of the $2n - i$ values in the union. Since the expected minimum of $\ell$ numbers drawn uniformly at random from $[0..N]$ is $N/(\ell + 1)$, the expected value of $m$ is
Improving the Sensitivity of MinHash Through Hash-Value Analysis

Figure 1 Probability distributions that two sets of size 100 share a common MinHash value, as a function of the size of their intersection (respectively 20, 50, and 100). The probability of small matching MinHash values are increase for relatively dissimilar sets.

\[ N/(2n - i + 1) \]

In the first case above, \( i = n \), so \( E[a] = N/(n + 1) \), while for the second case \( i = 1 \) and \( E[b] = N/(2n) \). Thus it is more likely that \( \min(a_1, a_2) \) corresponds to case (1) and \( \max(a_1, a_2) \) to case (2).

The situation is illustrated in Figure 1, which shows the probability of observing a given MinHash value \( a \) for three different intersection sizes. The probability of observing a MinHash value of 0 with a possible range \([0 \ldots 1000]\) is almost twice as high for two 100-element sets with a 20-element intersection than when the sets are identical. More similar pairs of sets, with larger intersection sizes, have greater probability of large matching MinHash values.

4 Expected Intersection Size as a Function of MinHash Value

In this section, we analyze the expected intersection size of two sets based on observing a particular matching MinHash value. For two \( n \)-element sets \( S \) and \( T \) where \( |S \cap T| = i \), \( J(S, T) = i/(2n - i) \). Thus analyzing the intersection size of \( S \) and \( T \) provides a result which can be alternately interpreted in the context of the Jaccard similarity of \( S \) and \( T \) for \( n \)-element sets.

Let \( S \) and \( T \) be two sets each of size \( n \). We limit our attention to the case where \( S \) and \( T \) are non-disjoint, which is necessary for MinHash values to legitimately collide, so \( S \cap T \neq \emptyset \). Further, we assume that range of \( h \) from \([0 \ldots N]\) is sufficiently large relative to \( n \) that we can discount the possibility of spurious collisions, namely that there does not exist \( s \in S \) and \( t \in T \) where \( h(s) = h(t) \) despite \( s \neq t \).

4.1 Prior Distribution

Determining the expected intersection size as a function of matching hash values requires knowledge of a prior distribution on the value of the intersection size. In the analysis below, we base our analysis on a uniform prior distribution, that all intersection sizes between \( S \) and \( T \) are equally likely. Thus for every \( i \in [1..n] \), \( \mathbb{P}[|S \cap T| = i] = 1/n \).
The uniform distribution appears most natural to us as a general prior, which is why we analyze this case below. That said, the true prior distribution differs with application, particularly as to whether pairs of randomly selected sets are likely to have large or small intersection sizes. The analysis below can be repeated for any particular well specified prior distribution in an analogous fashion.

### 4.2 Analysis

The probability of two sets sharing the MinHash value equals the Jaccard similarity index, that is

\[
P[M(S) = M(T) \mid i = |S \cap T|] = \frac{i}{2^i - i}.
\]

Because all intersections are equiprobable under our prior distribution, we have

\[
P[|S \cap T| = i \mid M(S) = M(T)] = \frac{i}{\sum_{j=1}^{n} 2^j - j}
\]

Note that given \(\ell\) random numbers \(x_1, \ldots, x_\ell\) uniformly drawn from \([1..N]\), for \(a \in [1..N]\), we have

\[
P[\min\{x_1, \ldots, x_\ell\} \leq a] = 1 - P[x_1 > a \& \ldots \& x_\ell > a] = 1 - \left(1 - \frac{a}{N}\right)^\ell.
\]

Then, the probability the MinHash is exactly \(a\) is given by

\[
P[\min\{x_1, \ldots, x_\ell\} = a] = \left(1 - \frac{a - 1}{N}\right)^\ell - \left(1 - \frac{a}{N}\right)^\ell.
\]

We now estimate the conditional probability \(P[|S \cap T| = i \mid M(S) = M(T) = a]\). We have

\[
P[|S \cap T| = i \mid M(S) = M(T) = a] = \frac{P[|S \cap T| = i] \wedge M(S) = M(T) \wedge (a = M(S \cup T))}{P[M(S) = M(T) \wedge (a = M(S \cup T))]} = \frac{P[|S \cap T| = i] \cdot P[a = M(S \cup T) \mid |S \cap T| = i]}{\sum_{j=1}^{n} P[M(S) = M(T) \wedge (a = M(S \cup T)) \mid |S \cap T| = i]}.
\]

The last rewrite follows because \(P[|S \cap T| = i] = 1/n\) is the same for all \(i\). To further simplify Eqn. 5, observe that

\[
P[M(S) = M(T) \mid |S \cap T| = i \wedge (a = M(S \cup T))] = P[M(S) = M(T) \mid |S \cap T| = i]
\]

because the event of sharing common MinHash \((M(S) = M(T))\) is independent of its value \((a)\) for a fixed intersection size. For the same reason, in the denominator,

\[
P[M(S) = M(T) \wedge (a = M(S \cup T)) \mid |S \cap T| = i] = P[M(S) = M(T) \mid |S \cap T| = i] \cdot P[a = M(S \cup T) \mid |S \cap T| = i]
\]

Eqn. 5 then rewrites to

\[
\frac{P[a = M(S \cup T) \mid |S \cap T| = i] \cdot P[M(S) = M(T) \mid |S \cap T| = i]}{\sum_{j=1}^{n} P[M(S) = M(T) \mid |S \cap T| = i] \cdot P[a = M(S \cup T) \mid |S \cap T| = i]}.
\]
Improving the Sensitivity of MinHash Through Hash-Value Analysis

Figure 2 The probability of intersection size of two sets of size 100 sharing a common MinHash value. The red curve shows the probability of having a given intersection size (formula (2)). The other curves show the same probability conditioned on the value $a$ of common MinHash (formula (7)), where the hash space is $[1..1000]$. Larger values of $a$ favor larger intersection sizes.

Using (4), (1), we obtain

$$\Pr[|S \cap T| = i \mid M(S) = M(T) = a] = \frac{i}{2^{n-i}} \left( \left( 1 - \frac{a-1}{N} \right)^{2n-i} - \left( 1 - \frac{a}{N} \right)^{2n-i} \right) \sum_{j=1}^{a} \frac{j}{2^{n-j}} \left( \left( 1 - \frac{a-1}{N} \right)^{2n-j} - \left( 1 - \frac{a}{N} \right)^{2n-j} \right)$$

(7)

Using (7), we can compute the expected intersection size as a function of the shared MinHash value:

$$\mathbb{E}[|S \cap T| \mid M(S) = M(T) = a] = \sum_{i=1}^{n} i \cdot \Pr[|S \cap T| = i \mid M(S) = M(T) = a]$$

(8)

As an illustration, Figure 2 shows probability distributions of intersection sizes without taking into account the common MinHash value (formula (2)) and knowing the common MinHash value (formula (7)). The figure demonstrates that larger common MinHash values provide an evidence for larger intersection sizes.

5 Hash Scoring for Sketch Similarity

In the classical MinHash scheme, the probability that two sets have matching MinHash is equal to the Jaccard similarity between the two sets. Thus, the fraction of matches taken over a number of trials provides an unbiased estimator of the Jaccard similarity. We have shown that the values of these matching MinHashes provides an orthogonal measure of similarity. The question is what the best way to combine these measures is.

We propose the following initial strategy. Traditional MinHash can be interpreted as averaging the values of 0/1 indicator variables, where a match of hash codes is represented by 1 and a mismatch by 0. We will replace the value associated with matching hashes by real values that over/underweight based on the value of shared MinHash. More specifically, a shared MinHash value will contribute with weight

$$\frac{\mathbb{E}[|S \cap T| \mid M(S) = M(T) = a]}{\mathbb{E}[|S \cap T|]}$$

(9)
Table 1 Improvement (in terms of the average reduction of absolute error) in estimating set intersection size by summing hash-weighted counts vs. equal weighting to estimate Jaccard similarity for different numbers of hash functions (rows) and set intersection sizes (columns). Bolded entries represent improvement over traditional MinHash estimation, representing $116/140 = 82.9\%$ of the non-trivial cells in the table.

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where the numerator is defined by Eqn 8, and

$$\mathbb{E}[|S \cap T|] = \sum_{i=1}^{n} P(|S \cap T| = i \mid M(S) = M(T))$$  \hspace{1cm} (10)$$

is the expected intersection size independent of MinHash value, i.e. implied by the prior distribution of intersection sizes.

5.1 Experimental Results

We performed a modest experiment to evaluate the performance of this technique, with the results reported in Table 1. We limited the experiment to small sets ($n = 20$), but consider a broad range of hash function counts ($1 \leq k \leq 20 = n$) and set similarities defined by intersection sizes from $1 \leq i \leq 20 = n$. Each cell represents the average difference in absolute error in estimating intersection size between MinHash with over/underweighting and traditional 0/1 counts, where each cell is averaged over 1,000 independent random trials.

We note that the rightmost column in Table 1 (intersection size 20 out of a possible 20) corresponds to identical sets, where the traditional Jaccard (and intersection size) estimate is always correct, leaving our proposed method with no room for possible improvement. But $116/140 = 82.9\%$ of the non-trivial cells show improvement over the traditional MinHash baseline.
Improving the Sensitivity of MinHash Through Hash-Value Analysis

Figure 3 The Kolmogorov-Smirnov test quantifies the difference between two probability distributions by the maximum $y$-distance gap between the two cumulative distribution functions. On the left, two samples from the same normal distribution. On the right, comparison of samples from uniform and normal distributions drawn over the same x-range.

6 Sketch Evaluation using the Kolmogorov-Smirnov Test

We now propose an alternate approach to improve the Jaccard similarity estimate offered by the classical MinHash approach, namely the fraction of matching MinHash values in $k$ trials. We seek to improve this estimate by analyzing the distribution of the values of the matching hashes from these trials to decide whether it is more likely to be over or under estimating the actual similarity. A key advantage of this approach over that of Section 4 is that does not require a prior distribution on the actual intersection sizes.

Our approach is based on the Kolmogorov-Smirnov (KS) statistical test [18, 23], which compares empirical cumulative distribution functions (CDFs) to assess whether two samples are drawn from the same underlying distribution. We will use it to compare the observed distribution of matching MinHash values against the theoretical distribution for the classical Jaccard estimate. The direction of the largest deviation suggests whether it is more likely an over or under estimate.

6.1 The Kolmogorov-Smirnov Test

In the KS-test, the empirical cumulative distribution function (CDFs) of the two different samples are plotted on the same chart. If the two samples are drawn from the same distribution, the ranges of $x$ values should largely overlap. An empirical CDF $\hat{F}(x)$ of a sample is defined as the fraction of the sample $\leq x$.

We seek to identify the value of $x$ for which the associated values of the two CDFs differ by as much as possible. The distance $D(\hat{F}, \hat{G})$ between two empirical CDFs $\hat{F}$ and $\hat{G}$ is the difference of the $y$ values at this critical $x$, formally stated as

$$D(\hat{F}, \hat{G}) = \max_x |\hat{F}(x) - \hat{G}(x)|$$

The more substantially that two samples differ in this fashion, the more likely it is that they were drawn from different distributions. Figure 3 (left) shows two independent samples from the same normal distribution. In contrast, Figure 3 (right) compares a sample drawn from a normal distribution against one drawn from the uniform distribution. The big gaps near the tails provide evidence that the two samples are drawn from different distributions.
The KS-test compares the value of \( D(\hat{F}, \hat{G}) \) against a particular target, declaring that two distributions differ at the significance level of \( \alpha \) when:

\[
D(\hat{F}, \hat{G}) > c(\alpha) \sqrt{\frac{n_1 + n_2}{n_1 n_2}}
\]

where \( c(\alpha) \) is a distribution-independent constant to look up in a table. In this paper, we use the ideas behind the KS-test for qualitative evaluation instead of precisely measuring statistical significance, and so will be interested in the direction of the deviation without this associated constant.

### 6.2 Application to MinHash Analysis

As explained above, the distribution of matching MinHash values differs as a function of the intersection size or (equivalently) Jaccard similarity between two sets of size \( n \). Recall that for \( \ell \) random numbers \( x_1, \ldots, x_\ell \) uniformly drawn from \([1..N]\), for \( a \in [1..N] \), we have

\[
F_\ell(x) = \mathbb{P}[\min\{x_1, \ldots, x_\ell\} \leq x] = 1 - \mathbb{P}[x_1 > x & \ldots & x_\ell > x] = 1 - \left(1 - \frac{x}{N}\right)^\ell \approx 1 - e^{-x\ell/N}. \tag{11}
\]

This defines the CDF on matching MinHash values. Comparing two sets \( A \) and \( B \), both of cardinality \( n \) with an intersection of size \( i \), any common MinHash value represents the smallest of \( \ell = 2n - i \) random values. Thus the distribution of matching MinHash values is defined by Eqn. 11, given an estimate for the union size \( \ell \). An important observation for us is that CDFs \( F_\ell \) are majorating one another, that is if \( \ell_1 > \ell_2 \), then \( F_{\ell_1}(x) > F_{\ell_2}(x) \) for any \( x \).

Estimates for the union size \( \ell \) and intersection size \( i \) follow from classical MinHash analysis. If \( m \) matching MinHash values are observed in \( k \) trials, \( m/k \) is an unbiased estimator of the Jaccard index \( \hat{i} = \frac{1}{2n-\ell} = \frac{2n-\ell}{k} \). Therefore, \( \ell \) and \( i \) are estimated respectively by

\[
\hat{i} = \frac{2nm}{k + m}, \quad \hat{\ell} = \frac{2nk}{k + m}.
\]

We can now employ the idea underlying the KS-test to evaluate how well the \( m \) observed MinHash values match the estimated distribution \( F_\ell(i) \). In doing that, we analyze the sign of the critical deviation

\[
D(F_\ell, \hat{F}) = F_\ell(\hat{x}) - \hat{F}(\hat{x}) \text{ for } \hat{x} = \text{argmax } |F_\ell(x) - \hat{F}(x)|,
\]

where \( \hat{F} \) is the empirical CDF obtained from the sample of matching MinHash values. When \( D \) is positive, this suggests that the regular MinHash estimate \( \hat{\ell} \) is an overestimate and therefore \( \hat{i} \) is an underestimate for the true intersection size. Conversely, a negative \( D \) provides an evidence that \( \hat{i} \) is an overestimate for the true intersection size. This reasoning is supported by the above-mentioned majorating property of CDFs, as it guarantees that the sign of \( D \) correctly defines whether the estimate \( \hat{\ell} \) should be increased or decreased for the KS-test statistic to be reduced and therefore for the estimated CDF to fit better the observed MinHash values. We thus propose the sign of \( D \) as a secondary signal to improve the accuracy of \( \hat{i} \) as an estimator for intersection size.

The running time of this test is \( O(m \log m) \) because we must sort the observed matching hash values to compute the CDF. It is only necessary to compare the distributions at the \( m \) sample points to identify the extremal points, with each comparison efficiently done using the exponential form of Eqn. 11. The magnitude of the deviation directly maps to a confidence value in the direction of change, with \( p \)-values obtainable using tables of \( c(\alpha) \) values from the standard KS-test. However, in the experiments below we propose to estimate correction direction from the sign of \( D \) independent of its magnitude.
Table 2 Performance (in terms of the fraction of correct direction predictions) of the KS-test-based over/under correction, as a function of the number of hash functions $k$ (shown in left column), and the true Jaccard similarity/intersection size (shown in first/second row). Generally speaking, the improvement is greatest at extreme values of similarity (either high or low), and with smaller numbers of hash functions.

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<td>.596</td>
<td>.584</td>
<td>.586</td>
<td>.550</td>
<td>.561</td>
<td>.591</td>
<td>.580</td>
<td>.608</td>
<td>.690</td>
<td>.660</td>
<td>.697</td>
<td>.610</td>
</tr>
</tbody>
</table>

6.3 Experimental Results

Table 2 summarizes the performance of our KS-based correction strategy over a wide range of hash counts (from $k = 1$ to $k = 1000$) and true Jaccard similarity (from 0.081 to 0.961). For each Jaccard similarity level, we constructed 10,000 pairs of 1000-element sets, each pair constructed to the appropriate level of similarity. We then constructed $k$ independent hash functions of these sets, and determined the number of matching MinHashes for these trials. We then performed the KS-test on the matching values to propose whether the actual Jaccard estimate should be higher or lower than the observed fraction of matches. We chose parameters of our tests (intersection size) so that to avoid the situation when the MinHash estimate exactly equals the true Jaccard similarity, making each case a fair binary trial.

Of the $14 \times 12 = 172$ entries in Table 2, 144 of them (83.7%) are greater than 0.5, meaning the adjustment breaks in the correct direction more often than not. The average accuracy ratio taken over all trials is 61.0%, substantially better than the baseline of 50%.

When employing large numbers of hash functions $k \geq 100$, our technique improves the estimate on average in 57 of 60 (95%) entries, and proves most beneficial in the middle regions where the Jaccard similarity is $\approx 0.5$. This is curious, because larger $k$ provides greater resolution on the fraction of matching hash values, thus reducing the quantization error of classical MinHash. But the KS-analysis also improves with more samples as $k$ increases, and continues to refine the similarity estimate even as $k = 1000$. Presumably in the limit as $k$ grows, the improvement over baseline will disappear, but it seems durable over the range of $k$ that appear in general applications.

That our best (and worst) performance occurs for very small $k$ reflects issues of quantization: for an intersection size of $n/2$ and $k = 5$, the best possible estimate must be wrong by at least 10%. As a statistical significance test, the KS-test was designed to be used with a meaningful number of samples per observed distribution. There are likely other statistical tests to do better with small (and maybe even large) values of $k$. 
7 Conclusions

We have demonstrated that the value of matching MinHash values provides additional information on the degree of similarity between pairs of sets. Our wins are small, but they are real. We believe that there exist better methods of integration to synthesize the mix of the number of matching hashes and their values into a more accurate measure of similarity and believe that this is a research direction worth pursuing. We note that even careful analysis of the values of the matching hash codes will be substantially less computationally expensive than that of obtaining the MinHash codes themselves, so these improvements will come at a little computational cost.

The MinHash values that do not match also contain some degree of signal concerning the similarity of two sets. Suppose the smallest hash values of two sets do not match, but are both unusually large, say a substantial fraction of the total range $N$. These large MinHash values signify that both sets exclude the same large fraction of possible elements from the universe, implying they must both be constructed from just a relatively small set of non-excluded elements. This conditioning increases the expected Jaccard similarity, despite the fact that the hash values do not match. We believe this signal to be very weak except in extreme cases, but its analysis may be part of a complete solution.

The theoretical success of MinHash depends strongly upon the elements of the sets being distinct. If multiplicity of elements should be taken into account, one should resort to the weighted variant of MinHash [26]. Extending our ideas to Weighted MinHash is another interesting direction of study for the future.

References


Suffix-Prefix Queries on a Dictionary

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Abstract

In the all-pairs suffix-prefix (APSP) problem, we are given a dictionary \( R \) of \( k \) strings, \( S_1, \ldots, S_k \), of total length \( n \), and we are asked to find the length \( SPL_{i,j} \) of the longest string that is both a suffix of \( S_i \) and a prefix of \( S_j \), for all \( i, j \in [1,k] \). APSP is a classic problem in string algorithms with many applications in bioinformatics. When all strings of the dictionary are over an integer alphabet of size \( \sigma \leq n^{O(1)} \), APSP can be solved in the optimal \( O(n + k^2) \) time with the use of the generalized suffix tree of the dictionary [Gusfield et al., Inf. Process. Lett. 1992].

In many bioinformatics applications, such as in sequence assembly, the size \( k \) of dictionary \( R \) is very large. In particular, \( k^2 \) usually dominates \( n \), and thus the \( k^2 \) factor is the bottleneck both in the time and in the space complexity of such applications. We thus initiate a holistic study on several data structure variants of APSP. In particular, we consider the following types of queries:

- **One-to-One** \((i,j)\): output \( SPL_{i,j} \).
- **One-to-All** \((i)\): output \( SPL_{i,j} \) for every \( j \in [1,k] \).
- **Report** \((i,\ell)\): output all distinct \( j \in [1,k] \) such that \( SPL_{i,j} \geq \ell \), where \( \ell \geq 0 \) is an integer.
- **Count** \((i,\ell)\): output the number of distinct \( j \in [1,k] \) such that \( SPL_{i,j} \geq \ell \), where \( \ell \geq 0 \) is an integer.
- **Top** \((i,K)\): output \( K \) distinct \( j \in [1,k] \) with the highest values of \( SPL_{i,j} \) breaking ties arbitrarily.

We assume the standard word RAM model of computation with word size \( w = \Omega(\log n) \) and an integer alphabet of size \( \sigma \leq n^{O(1)} \). We show the following upper bounds:

<table>
<thead>
<tr>
<th>Query</th>
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<tbody>
<tr>
<td>One-to-One ((i,j))</td>
<td>( O(n) )</td>
<td>( O(\log \log k) )</td>
<td>Theorem 11</td>
</tr>
<tr>
<td>One-to-All ((i))</td>
<td>( O(n) )</td>
<td>( O(k) )</td>
<td>Theorem 14</td>
</tr>
<tr>
<td>Report ((i,\ell))</td>
<td>( O(n) )</td>
<td>( O(\log n / \log \log n + \text{output}) )</td>
<td>Theorem 19(i)</td>
</tr>
<tr>
<td>Count ((i,\ell))</td>
<td>( O(n) )</td>
<td>( O(\log n / \log \log n) )</td>
<td>Theorem 19(ii)</td>
</tr>
<tr>
<td>Top ((i,K))</td>
<td>( O(n) )</td>
<td>( O(\log^2 n / \log \log n + K) )</td>
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We also present efficient algorithms for constructing these data structures.

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Introduction

The all-pairs suffix-prefix problem (APSP, in short) is a classic problem in string algorithms. APSP finds numerous applications in bioinformatics because it is the first step in sequence assembly [26, 37, 46, 8, 11]. Given a dictionary $R$ of $k$ strings, $S_1, \ldots, S_k$, of total length $n$, the APSP problem asks us to find, for each string $S_i$, $i \in [1, k]$, its longest suffix that is a prefix of string $S_j$, for all $j \neq i$, $j \in [1, k]$. Gusfield et al. [27] presented an algorithm running in the optimal $O(n(k^2))$ time for solving APSP, assuming all strings in $R$ are over an integer alphabet of size $\sigma \leq n^{O(1)}$. The algorithm is based on the generalization suffix tree [53] of $R$. Ohlebusch and Gog [39] gave another optimal algorithm which is based on the generalized suffix array [36] of $R$. Tustumi et al. [49] gave yet another optimal algorithm based on the generalized suffix array of $R$. Thus the common denominator of all existing optimal algorithms for APSP is that they rely on sorting the suffixes of all strings in $R$, and therefore they require space $\Omega(n)$ in any case and for any alphabet. In a very recent work, Loukides and Pissis [34] presented a different optimal algorithm, which is based on the Aho-Corasick automaton of $R$ [1], and it thus requires space linear in the size of the automaton.

Due to the practical relevance of APSP, there also exists a large body of works devoted to implementing algorithms for APSP that are suboptimal but practically fast on real-world datasets; see [25, 42, 33] and references therein for some of the state-of-the-art implementations. For a parallel implementation of the algorithm by Tustumi et al. see [35]. For approximate variants of APSP, under the Hamming or edit distance, see [44, 52, 32, 5, 47].

In many bioinformatics applications, such as in sequence assembly, the size $k$ of dictionary $R$ is very large. In particular, $k^2$ usually dominates $n$, and thus the $k^2$ factor is the bottleneck both in the time and the space complexity of such applications. For instance, in typical benchmark datasets\(^1\) for genome assembly using short DNA reads (fragments), $k$ is in the order of $10^6$ to $10^8$ and $n$ is in the order of $10^8$ to $10^{10}$. Hence $k^2$ dominates $n$ significantly.

We thus initiate a holistic study on several data structure variants of APSP. Let $\text{SPL}_{i,j}$ (short for suffix-prefix length), for any $i, j \in [1, k]$, denote the length of the longest string that is both a suffix of $S_i$ and a prefix of $S_j$. We consider the following types of queries:

- **One-to-One**$(i, j)$: output $\text{SPL}_{i,j}$.
- **One-to-All**$(i)$: output $\text{SPL}_{i,j}$ for every $j \in [1, k]$.
- **Report**$(i, \ell)$: output all distinct $j \in [1, k]$ such that $\text{SPL}_{i,j} \geq \ell$, where $\ell \geq 0$ is an integer.
- **Count**$(i, \ell)$: output the number of distinct $j \in [1, k]$ such that $\text{SPL}_{i,j} \geq \ell$, where $\ell \geq 0$ is an integer.
- **Top**$(i, K)$: output $K$ distinct $j \in [1, k]$ with the highest values of $\text{SPL}_{i,j}$ breaking ties arbitrarily.

By being able to answer different types of such queries efficiently, one may be able to design alternative algorithms, depending on the application in scope, which avoid the $k^2$ factor in their time or space complexity. Indeed, we stress that most works studying APSP from a practical perspective (e.g., [25, 42, 33]), in fact considered the $\ell$-APSP problem in their experimental part; namely, the problem in which we are asked to output only the $\text{SPL}_{i,j}$ values with $\text{SPL}_{i,j} \geq \ell$, for some integer $\ell \geq 0$, which, however, is given a priori and is fixed for all pairs $S_i, S_j$. This inflexibility would be surpassed should one have space-efficient (e.g., linear-space) data structures for answering these different types of queries fast.

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\(^1\) For example, see [http://gage.cbcb.umd.edu/data/index.html](http://gage.cbcb.umd.edu/data/index.html).
Our Results. We assume the standard word RAM model of computation with word size $w = \Omega(\log n)$ and an integer alphabet of size $\sigma \leq n^{O(1)}$. We show the following upper bounds:

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<td>$O(\log^2 n / \log \log n + K)$</td>
<td>Theorem 22</td>
</tr>
</tbody>
</table>

We also provide efficient construction algorithms for Theorems 11 and 14: Theorem 11 can be implemented in $O(n \log \log k)$ time and Theorem 14 can be implemented in $O(n)$ time. For Theorems 19 and 22, no guaranteed construction time is provided: the query times for Report, Count, and Top rely on the construction of a 2D rectangle stabbing data structure for reporting [45] and counting [28], but unfortunately the construction times for these data structures are not mentioned in [45] or [28]. However, by constructing the classic data structure for 2D rectangle stabbing [15], we obtain $O(n \log n)$ construction time, $O(n)$ words of space, $O(\log n + \text{output})$ query time for Report, $O(\log n)$ query time for Count, and $O(\log^2 n + K)$ query time for Top. We also make the following straightforward observation.

- Observation 1. The symmetric versions of One-to-All, Report, Count and Top, where we are given string $S_j$ as the query and we are asked to output information about SPL$_{i,j}$, for all $i \in [1, k]$, can be addressed by constructing the corresponding data structures for the dictionary $R'$ of $k$ strings $S'_1, \ldots, S'_k$, where $S' = S[S[1]] \cdots S'[2]S[1]$ denotes the reverse of string $S = S[1]S[2] \cdots S[|S|]$. Hence, the same space/query-time trade-offs can be achieved.

Related Work. In addition to the data structure variants of APSP that are studied here, two other versions of APSP have been studied in the literature. The first version consists in enumerating all pairwise suffix-prefix matches (not necessarily the longest ones) in decreasing order of their lengths. This version of the problem was solved by Ukkonen [50], who used this solution as the crux of his classic linear-time implementation of the greedy algorithm for constructing approximate shortest common superstrings. The second APSP version studied consists in enumerating the set of longest suffix-prefix matches (not however their association with the corresponding pairs of strings) [12]. Since any suffix-prefix match in this set is a prefix of some input string, the size of this set is $O(n)$. This version of the problem was solved in the optimal $O(n)$ time, independently, by Park et al. [40] and by Khan [29].

Although our work is inspired by real-world applications, the underlying data structure problems are also appealing from a theoretical perspective: (i) they are analogous to distance oracles for networks [48, 41, 17, 16, 13]; and (ii) they are special types of internal pattern matching (IPM) data structures [31, 30, 3, 14, 4]. For instance, an existing, more general, IPM data structure [30, 31] can be employed to answer One-to-One queries in $O(\log n)$ time using $O(n)$ words of space; see Section 2.3 for more details. By designing a specialized data structure for One-to-One, we obtain $O(\log \log k)$ query time using $O(n)$ words of space.

Paper Organization. In Section 2, we provide basic definitions and notation on strings. We also describe basic data structures for representing a dictionary, some more advanced data structures that are necessary to obtain our upper bounds, and a few previous solutions to APSP (variants). In Section 3, we provide the solution to One-to-One queries. In Section 4, we provide the solution to One-to-All queries. In Section 5, we provide the solutions to Report and Count queries. Finally, in Section 6, we provide the solution to Top queries.
2 Preliminaries

An alphabet $\Sigma$ is a finite nonempty set of $\sigma = |\Sigma|$ elements called letters. By $\Sigma^*$ we denote the set of all strings over $\Sigma$ including the empty string $\varepsilon$ of length 0. A string $S$ over $\Sigma$ is a sequence of letters of $\Sigma$. For a string $S = S[1] \cdots S[n]$ over $\Sigma$, by $n = |S|$ we denote its length. The fragment $S[i \ldots j]$ of $S$ is an occurrence of the underlying substring $P = S[i] \cdots S[j]$. We also say that $P$ occurs at (starting) position $i$ in $S$. A prefix of $S$ is a fragment of $S$ of the form $S[1 \ldots j]$ and a suffix of $S$ is a fragment of $S$ of the form $S[i \ldots n]$.

Let $M$ be a finite nonempty set of strings over $\Sigma$ of total length $m$. We call $M$ a dictionary. We define the trie of $M$, denoted by $\text{TR}(M)$, as a deterministic finite automaton that recognizes $M$. Its set of states (nodes) is the set of prefixes of the elements of $M$; the initial state (root node) is $\varepsilon$; the set of terminal states is $M$; and transitions (edges) are of the form $\delta(u, \alpha) = ua$, where $u$ and $ua$ are nodes and $\alpha \in \Sigma$. The size of $\text{TR}(M)$ is thus $O(m)$. The compacted trie of $M$, denoted by $\text{CT}(M)$, contains the root, the branching nodes, and the terminal nodes of $\text{TR}(M)$. The term compacted refers to the fact that $\text{CT}(M)$ reduces the number of nodes by replacing each maximal branchless path segment with a single edge, and that it uses a fragment of a string from $M$ to represent the label of this edge in $\Sigma$ words of space. The nodes of $\text{TR}(M)$ that are included in $\text{CT}(M)$ are called explicit; all other nodes are called implicit. The size of $\text{CT}(M)$ is thus $O(|M|)$. The most well-known form of compacted trie is the suffix tree described next.

2.1 Suffix Tree and Aho-Corasick Automaton

We are given a dictionary $R$ of $k$ strings, $S_1, S_2, \ldots, S_k$, whose total length is $n = |S_1| + |S_2| + \cdots + |S_k|$. Every string in $R$ is over an integer alphabet $\Sigma$ whose size $\sigma$ is polynomial in $n$, i.e., $\Sigma = \{1, 2, \ldots, n^{O(1)}\}$ and thus $\sigma \leq n^{O(1)}$. For constructing specialized data structures and answering internal pattern matching queries, non-trivial representations of $R$ (different than a simple set of strings) are usually more efficient.

Let us set $T_R := S_1 S_2 \cdots S_k$, where $S_1 < S_2 < \cdots < S_k$ are letters that are strictly lexicographically smaller than any letter from $\Sigma$ (and as such they do not belong to $\Sigma$).

Let $\text{ST}(S)$ denote the suffix tree of string $S$, that is the compacted trie of all the suffixes of $S$. For any node $v$ of $\text{ST}(S)$, by $\text{str}(v)$ we denote the concatenation of the edge labels on the path from the root to $v$, and by $d(v) = |\text{str}(v)|$ we denote the string depth of $v$. The suffix array $\text{SA}(S)$ of $S$ is the lexicographically sorted array of the set of suffixes of $S$, represented by their starting positions; see Figure 1 for an example.

Lemma 2 ([53, 22]). For any string $S$ of length $m$ over an integer alphabet of size $\sigma \leq m^{O(1)}$, the suffix tree and the suffix array of $S$ can be constructed in $O(m)$ time.

We also denote $\text{ST}_i = \text{ST}(S_i S_i)$ and $\text{ST}_R = \text{ST}(S_1 S_2 \cdots S_k)$; that is $\text{ST}_R$ is the generalized suffix tree [51] of the $k$ strings from $R$. The generalized suffix tree can be built in linear time; here, however, this more complicated construction is not needed since this compacted trie is equivalent to $\text{ST}(T_R)$ as the letters $S_i$ occur uniquely in this string (and hence a compacted edge containing any label $S_i$ must end at a leaf node).

Another useful representation of $R$ is given by its Aho-Corasick (AC) automaton [1]; the set of states of the AC automaton of $R$, denoted by $\text{AC}(R)$, corresponds to the set of the prefixes of the strings in $R$. Let $\text{node}(S)$ denote the node corresponding to string $S$. After reading an input string the automaton must be in a state corresponding to a suffix of this string (the longest one that is also a prefix of some string in $R$ and has a corresponding state); such a state always exists as $\varepsilon$ is always represented (recall $\varepsilon$ is the string of length 0). As such, the automaton $\text{AC}(R)$ is often represented by the trie $\text{TR}(R)$ with transitions $\delta(\text{node}(S), \alpha) =$
{node(Sα)} if Sα is a prefix of a string in R, and δ(node(S), ε) = {node(S′)}, where S′ is the longest suffix of S which is also a prefix of a string in R. The ε-transitions are called failure transitions. The existence of ε-transitions makes the automaton nondeterministic, and even though this nondeterminism can be avoided, we are going to actually employ those ε-transitions to construct the data structure for One-to-All queries.

Lemma 3 ([1], [20]). For any dictionary $R$ of $k$ strings of total length $n$ over an integer alphabet of size $σ \leq n^{O(1)}$, $AC(R)$ can be constructed in $O(n)$ time.

By $FT(R)$ we denote the so-called Failure Transition tree (FTtree) of $R$, introduced by Loukides and Pissis in [34] for solving the APSP problem: the FTtree nodes correspond to the states of the AC automaton (that is, to prefixes of strings in $R$), and the edges correspond to its ε-transitions with reversed direction. Notice that, since every state of $AC(R)$ has exactly one outgoing failure transition, $FT(R)$ is indeed a tree rooted at node($ε$). We additionally decorate every node $u$ of $FT(R)$ by a labeled interval $I_u = [i, j]_d$: $S_i, S_{i+1}, \ldots, S_j$ have as a common prefix the string of length $d$ represented by node $u$; see [34]. We will generally assume that $R$ is given lexicographically sorted at construction time; otherwise, the sorted version of $R$ can be produced in linear time using, for example, Lemma 3 or Lemma 2.

Example 4. Let $R = \{S_1, S_2, S_3, S_4\} = \{ACAA, ACAG, ACCG, CACA\}$ be a dictionary of $k = 4$ strings. The AC automaton and the FTtree of $R$ is shown in Figure 2. Consider the path from the root to leaf node $S_4$ (shown in red) in the FTtree of $R$, where the non-root nodes have the following labeled intervals $[i, j]_d$: $[1, 3], [4, 4], [1, 2], [4, 4]$. By recording the largest string depth $d$ of an interval containing $j$, for every $j \in [1, k]$, along this path, we compute all $SPL_{d,j}$: $SPL_{4,1} = 3$, $SPL_{4,2} = 3$, $SPL_{4,3} = 1$, and $SPL_{4,4} = 4$. Loukides and Pissis [34] showed how to compute this information, for all $i$, in $O(n + k^2)$ total time, thus solving the APSP problem optimally using only the FTtree of $R$.

2.2 Advanced Data Structures

Let $T$ be a rooted tree. A lowest common ancestor (LCA) query on $T$ for two given nodes $u$ and $v$, denoted by $w = LCA_T(u, v)$, returns the last (i.e., the lowest) common node $w$ on their paths from the root.

Lemma 5 ([9]). For any rooted tree $T$ with $m$ nodes, after $O(m)$-time preprocessing, we can answer $LCA_T$ queries in $O(1)$ time per query.
Figure 2 The AC automaton AC(R) (on the left) and FTtree FT(R) (on the right) of the dictionary of strings \(R = \{S_1, S_2, S_3, S_4\} = \{\text{ACAA, ACAG, AGGC, CACA}\}\). In AC(R), solid arrows correspond to transitions and dashed arrows to failure transitions. To avoid cluttering the figure, failure transitions to the start node in AC(R) have been omitted.

A rank and select data structure (also known as succinct indexable dictionary [43]) is a classic data structure, constructed over an array \(A\) of length \(m\) over alphabet \([1, \sigma]\), which supports two types of queries:

- \(\text{rank}_A(i, x) = |\{\ell \in [1, x] : A[\ell] = i\}|\), for \(i \in [1, \sigma]\) and \(x \in [1, m]\);
- \(\text{select}_A(i, x) = \min\{\ell \in [1, m] : \text{rank}_A(i, \ell) = x\}\), for \(i \in [1, \sigma]\) and \(x \in [1, m]\).

In other words, \(\text{rank}_A(i, x)\) returns the number of elements with value equal to \(i\) occurring at positions in \([1, x]\) of \(S\), while \(\text{select}_A(i, x)\) returns the position of the \(x\)th element of \(A\) with value equal to \(i\).

**Lemma 6 ([7, 38, 18]).** For any array \(A = A[1 \ldots m]\) over \([1, \sigma]\), \(\sigma \leq m\), after \(O(m \log \log \sigma)\)-time preprocessing, we can construct a data structure of \(O(m)\) words of space that supports \(O(\log \log \sigma)\)-time rank and select queries on \(A\).

Let \(T\) be a rooted tree of \(m\) nodes with integer weights on nodes. Further assume that the weight of every node of \(T\) satisfies the min-heap property: the weight of each node is greater than or equal to the value of its parent (the smallest weight is hence at the root). A weighted ancestor (WA) query for a given node \(u\) of \(T\) and an integer \(d\), denoted by \(w = \text{WA}_T(u, d)\), returns its deepest ancestor \(w\) whose weight is at most \(d\) [23]. This problem is the generalization of the classic predecessor search problem on rooted trees. In the special case when \(T\) is a suffix tree and the nodes are weighted by string depth, the problem admits an optimal solution due to the recent result of Belazzougui et al. [6] (see also [24]).

**Lemma 7 ([6]).** For any suffix tree \(T\) with \(m\) nodes weighted by string depth, after \(O(m)\)-time preprocessing, we can answer \(\text{WA}_T\) queries in \(O(1)\) time per query.

In this special case, the ancestor at string depth exactly \(d\) may be an implicit node of \(T\), in which case the query outputs its closest explicit ancestor.

### 2.3 Previous Solutions

**\(O(n + k^2)\)-time Algorithm for APSP.** We describe the optimal solution to APSP given by Gusfield et al. in [27]. We set \(T_R := S_1S_2S_3\ldots S_k\), where \(S_1 < S_2 < \cdots < S_k\) are letters that are strictly lexicographically smaller than any letter from \(\Sigma\). We start by constructing the suffix tree \(\text{ST}_R = \text{ST}(T_R)\). Using a DFS traversal on \(\text{ST}_R\), we construct lists \(L(v)\) for all nodes \(v\) of \(\text{ST}_R\); \(L(v)\) stores all \(i\) such that the suffix of length \(d(v)\) of string \(S_i\) is \(\text{str}(v)\). Consider a string \(S_j\) from \(R\) and focus on the path \(P_j\) from the root of \(\text{ST}_R\) to the leaf node representing the longest suffix of \(S_j\), i.e., the entire string \(S_j\). Let \(v\) be a node on \(P_j\). A suffix of string \(S_i\) of length \(d(v)\) is a prefix of string \(S_j\) of the same length if and only if \(i\) is
in $L(v)$. However, for each index $i$, we want to record the deepest node $v$ on $P_j$ such that $i$ is in $L(v)$. It then follows that $d(v) = \text{SPL}_{i,j}$. In order to achieve a linear-time complexity, we perform another DFS maintaining $k$ stacks (one for each $S_i$). Upon visiting $v$, we push it on stack $i$ for every $i \in L(v)$. When the leaf node representing the entire string $S_j$ is reached, we scan the $k$ stacks and record, for each index $i$, the current top of the $i$th stack. When $v$ is reached in a backward edge traversal, we pop the top of any stack whose index is in $L(v)$. We obtain the following result.

- **Lemma 8 ([27]).** For any dictionary of $k$ strings of total length $n$ over an integer alphabet of size $\sigma \leq n^{O(1)}$, $\text{APSP}$ can be solved in the optimal $O(n + k^2)$ time.

In what follows, we assume that $k \geq \sqrt{n}$; otherwise, when $k < \sqrt{n}$, Lemma 8 implies an optimal solution to our data structure problems (linear preprocessing time, linear size and time-optimal queries), which precomputes and stores all answers.

**Internal Prefix-Suffix Queries for One-to-One.** Kociumaka considered the following data structure problem in [30]: Given two fragments $x$ and $y$ of a string $T$ and a positive integer $d$, report all suffixes of $y$ of length between $d$ and $2d - 1$ that also occur as prefixes of $x$ (represented as an arithmetic progression of their lengths). This is the Internal Prefix-Suffix Queries problem. Kociumaka showed the following result (see also [31]).

- **Lemma 9 (Theorem 1.1.3 in [30]).** For any string $T$ of length $m$ over an integer alphabet of size $\sigma \leq m^{O(1)}$, after $O(m)$-time preprocessing, we can answer Internal Prefix-Suffix Queries in $O(1)$ time per query.

By employing Lemma 9 on $T_R$, after an $O(n)$-time preprocessing, we can answer One-to-One queries in $O(\log(\min(|S_i|, |S_j|))) = O(\log n)$ time. In particular, we query for $x = S_j$, $y = S_i$, and $d = 2^\ell$, for all integers $0 \leq \ell \leq \log \min(|S_i|, |S_j|)$, to compute a representation of all the suffixes of $S_i$ that are also prefixes of $S_j$ and then return the length of the longest one as $\text{SPL}_{i,j}$. We obtain the following result, which we improve in Section 3.

- **Corollary 10.** For any dictionary of $k$ strings of total length $n$ over an integer alphabet of size $\sigma \leq n^{O(1)}$, we can construct a data structure of $O(n)$ words of space answering One-to-One queries in $O(\log n)$ time.

### 3 Answering One-to-One Queries

**Main Idea.** Say we want to find the longest suffix of $S_i$ that is a prefix of $S_j$. We first find the maximal longest common prefix between $S_j$ and any suffix of $S_i$. Say this suffix is $S_i[q..|S_i|]$ and we have that $S_i[q..q + r - 1] = S_j[1..r]$ is this longest common prefix. If this prefix is the whole $S_i[q..|S_i|]$, i.e., $|S_i| = q + r - 1$, then $r$ is clearly the answer. If this longest common prefix is not a suffix of $S_i$, i.e., $|S_i| > q + r - 1$, then the answer is the longest prefix of $S_i[q..q + r - 1]$, that is also a suffix of $S_i$.

Recall that $ST_i = ST(S_i, \$)$ and $ST_R = ST(T_R)$. Consider the path in $ST_R$ obtained by reading $S_j$ from its root (this path ends in a leaf node). When spelling any suffix of $S_i$ that is also a prefix of $S_j$ in $ST_R$ we use exactly the same path and end by going out of it when reading $\$$. This means, that $\text{SPL}_{i,j}$ is represented by the lowest node on this path that has an outgoing edge with label $\$$. In the following we focus on enhancing $ST_R$ and $ST_i$, for all $i \in [1, k]$, to obtain a data structure that allows finding the string depth of such a node (equal to $\text{SPL}_{i,j}$) efficiently. We will prove the following result.
Theorem 11. For any dictionary of $k$ strings of total length $n$ over an integer alphabet of size $\sigma \leq n^{O(1)}$, we can construct a data structure of $O(n)$ words of space answering One-to-One queries in $O(\log \log k)$ time. The data structure can be constructed in $O(n \log \log k)$ time.

Let us start with a straightforward auxiliary lemma.

Lemma 12. For any dictionary of $k$ strings $S_1, \ldots, S_k$ of total length $n$ over an integer alphabet of size $\sigma \leq n^{O(1)}$, in $O(n)$ time we can construct a data structure of $O(k)$ words of space that answers queries of the type “Is $S_j$ a suffix of $S_i$?” in $O(1)$ time.

Proof. Let $X^r$ denote the reverse of string $X$, i.e., $X^r = X[|X|] \cdots X[1]$. We first sort $S_1^r, \ldots, S_k^r$ lexicographically, and store for each $j \in [1, k]$ a value $\text{rlex}[j] \in [1, k]$ equal to the rank of $S_j^r$ in this sorted list. $S_j$ is a suffix of $S_i$ if and only if $S_j^r$ is a prefix of $S_i^r$. The crucial property of this ordering is that all the strings such that $S_j^r$ is their prefix form an interval from the position $\text{rlex}[j]$ to a position $\text{rlex}[j] + |S_j^r| - 1$, where $|S_j^r|$ is the total number of strings $S_1^r, \ldots, S_k^r$ starting with $S_j^r$; that is, $\text{rlex}[j] + |S_j^r|$ is the position of the first string having a longest common prefix with $S_j^r$ shorter than $|S_j^r|$. The values $\text{rlex}[j]$ and $|S_j^r|$, for all $j \in [1, k]$, can be computed in $O(n)$ time [19].

As for the querying, for any $i, j$, we have that $S_j$ is a suffix of $S_i$ if and only if $\text{rlex}[j] \leq \text{rlex}[i] < \text{rlex}[j] + |S_j^r|$, which is checked in $O(1)$ time. The total size of arrays $l$ and $\text{rlex}$ is $\Theta(k)$.

Construction. We start the construction of the data structure by constructing the data structure underlying Lemma 12. We also construct $\mathbf{ST}_R$ and $\mathbf{ST}_i$, for all $i \in [1, k]$, using Lemma 2. We enhance $\mathbf{ST}_R$ with the data structure for LCA queries underlying Lemma 5, and link the leaf nodes originating from suffixes of $S_i$ with the corresponding leaf nodes of $\mathbf{ST}_i$, for all $i \in [1, k]$. We construct an array $A = A[1 \ldots |T_R|]$ over $[1, k]$ such that $A[i] = i$ if the $i$th leaf node (from the left) of $\mathbf{ST}_R$ originates from a suffix of $S_i$; since the leaf nodes are ordered according to the lexicographic order of the suffixes they originate from, array $A$ can be easily extracted from $\mathbf{SA}(T_R)$ constructed by means of Lemma 2. We enhance array $A$ with the rank and select data structure underlying Lemma 6. We link the leaf nodes of $\mathbf{ST}_R$ with the corresponding elements of $A$. For each $\mathbf{ST}_i$, we construct the data structure for WA queries underlying Lemma 7. For every node $v$ of $\mathbf{ST}_i$, we store the string depth of its closest ancestor (including $w$ itself) that has an outgoing edge with label $\$i$ and hence corresponds to a suffix of $S_i$; since the root always has such an edge, this assignment is always well-defined. In order to efficiently compute and store all those values, we simply process the information through the tree in a top-down manner. This completes the construction.

The part of the data structure that relies on Lemmas 2, 5, 7, and 12 is implemented in $O(n)$ time and it occupies $O(n)$ words of space. By Lemma 6, array $A$ occupies $O(n)$ words of space, and it can be implemented in $O(n \log \log k)$ time as it stores $k$ distinct values.

Querying. Consider a One-to-One $(i, j)$ query; that is, we want to compute $\text{SPL}_{i,j}$, the length of the longest suffix of $S_i$ that is a prefix of $S_j$. Let $x$ be the position in array $A$ that corresponds to the leaf node $l_j$ of $\mathbf{ST}_R$ reached after conceptually reading $S_j$. We first check if the entire $S_j$ is a suffix of $S_i$ by means of Lemma 12. If this is the case then we return $\text{SPL}_{i,j} = |S_j|$. If this is not the case (inspect Figure 3), we perform the following sequence of queries, $\text{select}_A(i, \text{rank}_A(i, x))$, which finds the position $y$ in array $A$ that corresponds to the leaf node $r_j$; this corresponds to the suffix of $S_j$ that is closest to the left of $l_j$. We then compute the lowest common ancestor of $r_i$ and $l_j$: $v_{i,j} = \text{LCA}_{\mathbf{ST}_A}(r_i, l_j)$. If node $v_{i,j}$ has an outgoing edge labeled with $\$i$, which ends at $r_i$, then we return $\text{SPL}_{i,j} = d(v_{i,j})$ (this
Figure 3 An illustration of the One-to-One \((i,j)\) query algorithm. The node \(v_{i,j}\), which is explicit in \(ST_R\) but implicit in \(ST_i\), has an outgoing edge labeled with $i$, and hence the string depth \(d(v_{i,j})\) of node \(v_{i,j}\) is the answer to the query.

Figure 4 An illustration of the One-to-One \((i,j)\) query algorithm. The closest ancestor of node \(v_{i,j}\), which is explicit in \(ST_R\) but implicit in \(ST_i\), with an outgoing edge labeled with $i$ is node \(u\) and hence the string depth \(d(u)\) of node \(u\) is the answer to the query.

is the case in Figure 3). We check this by checking whether \(d(r_i) = d(v_{i,j}) + 1\). If \(v_{i,j}\) does not have such an outgoing edge (this is the case in Figure 4), we locate the explicit node corresponding to \(v_{i,j}\) in \(ST_i\) (or its closest explicit ancestor if it is implicit) by asking a WA query: \(w = \text{WA}_{ST_i}(r_i, d(v_{i,j}))\). Finally, we return the string depth of the closest ancestor of \(w\) with an outgoing edge labeled $i$ as \(\text{SPL}_{i,j}\); recall that every node of \(ST_i\) stores this information.

The time complexity of the query is \(O(\log \log k)\); the bottleneck is the complexity of the rank and select queries on \(A\) – all other operations take constant time. Let us now explain why the faster \(O(1)\)-time select and \(O(1 + \log \log \log n)\)-time rank queries presented in [7], where \(w\) is the machine word, cannot improve our query time further. The size of the problem is \(\Theta(n)\), hence the size of the machine word in the word-RAM model is \(\Theta(\log n)\), thus the query time equals \(O(1 + \log \log \log n)\). However, we have assumed that \(k \geq \sqrt{n}\) (otherwise the structure of Lemma 8 implies an optimal solution – linear size and constant time queries for the One-to-One queries), hence this is equal to \(O(1 + \log \log k) = O(\log \log k)\) as stated.

Correctness. Recall that the answer to One-to-One \((i,j)\) equals to the string depth of the closest ancestor of \(l_j\) in \(ST_R\) that has an outgoing edge labeled with $i$. By construction, this ancestor ends on the right of \(l_j\) only if the entire \(S_j\) is a suffix of \(S_i\), which we check separately. Otherwise, this ancestor is also an ancestor of \(r_i\) (which is on the left of \(l_i\)) as $i$
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goes out of the path from the root to \(l_j\) to the left (by construction, it is lexicographically smaller than the next letter on this path), and hence this edge labeled with \(\$\) must end either in \(r_i\) or further to the left (by the definition of \(r_i\)). As an ancestor of \(l_j\) and \(r_i\), it is also the closest ancestor of \(v_{i,j}\) with such an outgoing edge; the latter actually exists (possibly as an implicit node) in \(ST_i\) (unlike \(l_j\)). The final steps of the query algorithm find the string depth of the node corresponding to the searched ancestor in \(ST_i\) (string depth is a shared property of the corresponding nodes).

We have arrived at Theorem 11. Note that the construction time for our data structure is \(O(n \log \log k)\). The bottleneck for the construction time is the construction time for the rank and select data structure (Lemma 6).

4 Answering One-to-All Queries

The spine of the data structure described in this section is \(FT(R)\), the FTtree of \(R\) (see Section 2). Recall that for each node in \(FT(R)\) (representing each prefix of a string \(S_i\)), we store information about which strings from \(R\) it is a prefix of (see Figure 2).

Main Idea. The Aho-Corasick lemma [1] states that for any two nodes, \(\text{node}(U)\) and \(\text{node}(V)\), in \(AC(R)\), we have a failure transition from \(\text{node}(U)\) to \(\text{node}(V)\) if and only if \(V\) is the longest suffix of \(U\) that is also a prefix of some string in \(R\). As a consequence, in \(FT(R)\), \(\text{node}(S)\) is an ancestor of \(\text{node}(S')\) if and only if \(S\) is a suffix of \(S'\) (and both are prefixes of some strings from \(R\) as nodes of \(FT(R)\)). Thus the path from \(\text{node}(\epsilon)\) (the root) to \(\text{node}(S_i)\) in \(FT(R)\) contains exactly the nodes \(\text{node}(S)\) such that \(S\) is a suffix of \(S_i\) and a prefix of some string in \(R\). Those nodes are ordered according to the string length, hence the nodes closer to \(\text{node}(S_i)\) on this path will correspond to longer suffix-prefix matches.

A One-to-All(i) query can thus be answered by simply reading the path from the root to \(\text{node}(S_i)\) recording, for each \(j \in [1, k]\), the last node on the path corresponding to a prefix of \(S_j\). The space occupied by \(FT(R)\) is in \(O(n)\); and such a query algorithm can take \(\Theta(|S_i|)\), that is even \(\Theta(n)\) time. Hence, by such an algorithm, we would not really gain anything from constructing \(FT(R)\) in the preprocessing. On the other extreme, by running this algorithm not for a single path, but for the whole \(FT(R)\) using a DFS traversal, we can precompute the answers for all the values of \(i \in [1, k]\) in \(O(n + k^2)\) total time (and space), and then answer a query in \(O(k)\) time by simply outputting the \(k\) stored values; this would not be faster than using the algorithm by Gusfield et al. [27] or the one by Loukides and Pissis [34]. We will augment \(FT(R)\) to obtain a more efficient solution combining the space efficiency of the first approach with the low query time of the second one.

A \(\tau\)-micro-macro decomposition, introduced for rooted binary trees in [2], and then generalized for rooted general trees in [10] (after an appropriate mapping), is a partition of a rooted tree \(T\) of \(N\) nodes into \(O(N/\tau)\) connected subtrees, called micro trees. In the case of binary trees each micro tree is of size at most \(\tau\) and at most two of its nodes are adjacent to nodes in other micro trees. These nodes are referred to as top and bottom boundary nodes of the micro tree. The top boundary node is chosen as the root of the micro tree. The macro tree is a rooted tree of size \(O(N/\tau)\) whose nodes correspond to micro trees as follows (inspect Figure 5): The top boundary node \(t(C)\) of a micro tree \(C\) is connected to a boundary node \(\text{parent}(C)\) in the parent micro tree (apart from the root). The boundary node \(t(C)\) might also be connected to a top boundary node of a child micro tree, which we denote by \(\text{child}(C)\). Such a \(\tau\)-micro-macro decomposition can be computed in \(O(N)\) time for binary [2] and general [10] rooted trees. We summarize the above discussion in the lemma below.
Lemma 13 ([2, 10]). For any rooted tree $T$ with $N$ nodes and for any integer $\tau \in [1, N]$, the $\tau$-micro-macro decomposition of $T$ can be computed in $O(N)$ time.

We will prove the following result.

Theorem 14. For any dictionary of $k$ strings of total length $n$ over an integer alphabet of size $\sigma \leq n^{O(1)}$, we can construct a data structure of $O(n)$ words of space answering One-to-All($i$) queries in $O(k)$ time. The data structure can be constructed in $O(n)$ time.

Construction. We start the construction of the data structure by constructing $\FT(R)$ from $\AC(R)$ using Lemma 3. We compute the $\tau$-micro-macro decomposition of $\FT(R)$, for a parameter $\tau$ defined later, using Lemma 13. For each node $u$ of the $\FT(R)$, corresponding to a prefix $S$ of some string $S_i$ in $R$, we store the labeled interval $I_u$. For each boundary node in the $\tau$-micro-macro decomposition of $\FT(R)$, we store an array of $k$ integers, which for each $i \in [1, k]$, stores the string depth of its lowest ancestor node($S$) such that $S$ is a prefix of $S_i$. The additional size for storing this information in all the boundary nodes is $O(k \cdot n/\tau)$. We compute these arrays by performing a DFS over $\FT(R)$ with a set of $k$ stacks, one for every string in $R$, storing the string depths of ancestors of the visited node of each type (which $S_i$ they originate from). As there are only $2n$ updates of the stacks (each prefix of a string $S_i$ is stored and removed once from the $i$th stack) and the information is stored by simply reading the top values of the $k$ stacks, the total computation time is bounded by $O(n + k \cdot n/\tau)$.

Querying. Let us start with the following observation from [34] (inspect also Figure 2).

Observation 15 ([34]). Let $u$ and $v$ be two non-root nodes of $\FT(R)$ with labeled intervals $I_u = [i_u, j_u]_{d(u)}$ and $I_v = [i_v, j_v]_{d(v)}$, respectively, and such that $u$ is an ancestor of $v$. Then $d(u) < d(v)$ and either $[i_u, j_u]$ contains $[i_v, j_v]$ or $[i_u, j_u]$ and $[i_v, j_v]$ do not intersect.

Consider a One-to-All($i$) query; that is, we want to compute an array of length $k$, which stores $\SPL_{i,j}$, for all $j \in [1, k]$. We start by finding the closest boundary node on the path from the root to $\node(S_i)$; that is, the top boundary node of the micro tree containing $\node(S_i)$. On the path between this top boundary node and $\node(S_i)$, there are at most $\tau$ nodes. We compute the information coming from just those nodes in $O(k + \tau)$ time with a
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sweep line approach: there are \( \mathcal{O}(\tau) \) (labeled) intervals from \([1, k]\), the intervals are labeled by different values (string depth), but, by Observation 15, two intervals are either disjoint or the one with the larger string depth is contained in the one with the smaller one. Thus, it is enough to hold the active intervals on a stack to keep track of the longest possible suffix-prefix match: the interval on the top of the stack has the highest value and will end the soonest. The solution is then obtained as the position-wise maximum of the computed array and the array stored in the top boundary node, which we compute in \( \mathcal{O}(k) \) time.

**Correctness.** The correctness of the algorithm follows by the Aho-Corasick lemma (see also the discussion of the “main idea” paragraph above).

The data structure occupies \( \mathcal{O}(n + k \cdot n/\tau) \) words of space and supports One-to-All queries in \( \mathcal{O}(k + \tau) \) time. By setting \( \tau \) to \( k \) (or to \( ck \), for some positive constant \( c \) that balances the operation costs more efficiently) we obtain the complexities claimed in Theorem 14. Note that the data structure is constructed in \( \mathcal{O}(n + k \cdot n/\tau) \) time, which is \( \mathcal{O}(n) \) for \( \tau = \Theta(k) \). Thus the presented data structure for One-to-All queries is optimal.

5 Answering Report and Count Queries

In this section we are going to use \( \text{ST}_R \) again. This time, however, instead of augmenting \( \text{ST}_R \) with an LCA data structure and linking its nodes with the rank and select array, we are going to link the nodes with rectangles and employ classic results from computational geometry for reporting (see Lemma 16) and counting (see Lemma 17).

Let \([x_1, x_2] \times [y_1, y_2]\) denote a rectangle in a 2D space with edges parallel to the axes, where the intervals \([x_1, x_2]\) and \([y_1, y_2]\) are the projections of this rectangle to the \(x\)-axis and \(y\)-axis, respectively. In the reporting version of the 2D rectangle stabbing problem [15], we are given a set \( S \) of \( n \) rectangles to preprocess, so that when we are given a query point \( q = (x, y) \), we report the subset \( Q \subseteq S \) of rectangles \([x_1, x_2] \times [y_1, y_2]\) that contain \( q \): \( x_1 \leq x \leq x_2 \) and \( y_1 \leq y \leq y_2 \). In the counting version of 2D rectangle stabbing, we are asked to return \(|Q|\).

▸ **Lemma 16** ([45]). For any set \( S \) of \( n \) rectangles, we can construct a data structure of \( \mathcal{O}(n) \) words of space answering 2D rectangle stabbing reporting queries in \( \mathcal{O}(\log n / \log \log n + f) \) time, where \( f \) is the output size \(|Q|\).

2D rectangle stabbing counting is known to be reducible to 2D orthogonal range counting [21], and such a data structure for 2D orthogonal range counting can be found in [28].

▸ **Lemma 17** ([21, 28]). For any set \( S \) of \( n \) rectangles, we can construct a data structure of \( \mathcal{O}(n) \) words of space answering 2D rectangle stabbing counting queries in \( \mathcal{O}(\log n / \log \log n) \) time.

**Main Idea.** For every suffix \( S \) of a string in \( R \) that is represented by a node in \( \text{ST}_R \), we define a rectangle in 2D space: the \( x \) dimension corresponds to the lexicographically sorted list of all suffixes of strings in \( R \) whose prefix is \( S \); and the \( y \) dimension corresponds to interval \([0, |S|]\). A Report (resp. a Count) query is defined by two parameters, which form a point in the 2D space: \( i \) corresponds to string \( S_i \) in the same sorted list (\( x \) dimension) and \( \ell \) corresponds to the smallest length of interest (\( y \) dimension). By reporting (resp. counting) all rectangles enclosing this point (Lemmas 16 and 17), we locate all suffix-prefix matches. Extra care, however, needs to be taken in order to avoid double reporting (resp. counting).
Construction. We start the construction of the data structure by constructing $ST_R$ using Lemma 2. Let $u$ be an explicit or implicit node of $ST_R$ that is the parent of a leaf node reached with $S_i$: the labels of the path from root to $u$ form a suffix of $S_i$. For every such node $u$ and every $i$, we create a tuple $(L(u), R(u), d(u), i)$, where $L(u)$ and $R(u)$ are the (pre-order rank of) the leftmost and the rightmost leaf node under $u$, respectively. Note that such a node may correspond to multiple tuples for different $i$ values – this occurs when distinct elements of $R$ share the same suffix. There are exactly $n$ such tuples (one for every suffix) coming from $ST_R$ and we can compute them in $O(n)$ total time using a DFS traversal.

Recall that if we spell $S_j$ in $ST_R$ and the obtained leaf node $v$ has an ancestor of string depth $\ell$ which has an outgoing edge with label $S_i$, then $SPL_{i,j} \geq \ell$. The same property ($SPL_{i,j} \geq \ell$) can be expressed by $L(v) \in [L(u), R(u)]$ (namely, $u$ is an ancestor of $v$), and $\ell \in [0, d(u)]$ (namely, the string depth of $u$ is at least $\ell$) for a tuple $(L(u), R(u), d(u), i)$. Now note that $(L(u), R(u), d(u), i)$ forms a rectangle, whose identifier is $i$. In particular, $(L(u), R(u), d(u), i)$ can be viewed as rectangle $[L(u), R(u)] \times [0, d(u)]$ with satellite data $i$.

Now consider constructing the 2D rectangle stabbing data structure for reporting (resp. counting) for these $n$ rectangles, and then ask the query for a point $(L(v), \ell)$, where $v$ is the leaf node reached from the root by conceptually reading $S_j$. The data structure will report (resp. count) all of the suffixes of $S_i$, for $i \in [1,k]$, of length at least $\ell$ that are also prefixes of $S_j$. Unfortunately, such a solution differs from the expected results of $\text{Report}(i, \ell)$ and $\text{Count}(i, \ell)$ in the following two ways:

1. Instead of finding all $j \in [1,k]$ such that $SPL_{i,j} \geq \ell$ for a given $i$, we find all such $i \in [1,k]$ for a given $j$. This issue is addressed by Observation 1, which states that $\text{Report}(i, \ell)$ and $\text{Count}(i, \ell)$ reduce trivially to the problems considered here, denoted by $\text{Report}'(i, \ell)$ and $\text{Count}'(i, \ell)$, respectively (recall that the $r$ superscript refers to reversing the input strings);

2. If there are multiple prefixes of $S_j$ of length at least $\ell$ that are also suffixes of $S_i$, then we will report (resp. count) each of them leading to double reporting (resp. counting).

Although one may actually be interested in reporting or counting those multiple suffix-prefixes, in this paper, we are only interested in the longest ones. We address this issue by modifying the rectangles before the construction.

As mentioned earlier the first issue is resolved by Observation 1. To solve the second issue, we have to make the set of rectangles, for a single $i \in [1,k]$, pairwise disjoint while leaving their union unchanged. Notice that two such non-disjoint rectangles must come from a pair of nodes $u$ and $w$ in an ancestor-descendant relationship. An easy solution is to take, for every node $w$ which has an outgoing edge with label $S_i$, its closest ancestor $u$ which also has an outgoing edge with label $S_i$, and change the $[L(w), R(w)] \times [0, d(w)]$ rectangle into $[L(w), R(w)] \times [d(u) + 1, d(w)]$; inspect Figure 6. Since the part $[L(w), R(w)] \times [0, d(u)]$ is already contained in $[L(u), R(u)] \times [0, d(u)]$ the union remains unchanged, and since $u$ is the closest such ancestor, the other rectangles (for this $i$) cannot have a nonempty intersection with the newly obtained one (the intersection with the ones coming from the descendants of $w$ is empty after the modification of those rectangles). We can perform these modifications with a single DFS traversal with $k$ stacks of nodes from the path from the root to the currently processed node, which has an outgoing edge with label $S_i$, $i \in [1,k]$. A more complicated solution is obtained by replacing the two rectangles $[L(u), R(u)] \times [0, d(u)]$ and $[L(w), R(w)] \times [0, d(w)]$ with three rectangles: $[L(u), L(w) - 1] \times [0, d(u)]$, $[L(w), R(w)] \times [0, d(w)]$. 

\footnote{[L(u), R(u)] is also known as the suffix array interval of node u.}
Two intersecting rectangles implied by \( ST_R \).

(a) Two intersecting rectangles implied by \( ST_R \).

(b) Two ways to make the rectangles disjoint.

**Figure 6** On the bottom left part, the rectangles obtained from two nodes \( u \) and \( w \) of \( ST_R \) (top left), both having an outgoing edge with label \( S_i \), forming a suffix-prefix match of \( S_i \) and \( S_j \) for node \( v \) reached by reading \( S_j S_i \) from the root. The rectangles have a nonempty intersection. To avoid double reporting (or double counting), we make the rectangles disjoint while leaving their union unchanged. We can do this (by taking the intersection once) in two ways (on the right): a simple one (top) or a more complicated one (bottom), which allows us to efficiently output \( SPL_{i,j} \).

\[ [0, d(w)] \] and \([R(w) + 1, R(u)] \times [0, d(u)]\]; inspect Figure 6. Unlike the previous construction, a single rectangle can be spliced into smaller ones many times (a node can be a direct ancestor of many other nodes); at the same time a single rectangle can splice only its direct ancestor, hence the number of rectangles obtained this way is bounded from above by \( 2n \). This set of modified intervals can be obtained similarly: in a DFS traversal, when a node which has an outgoing edge with label \( S_i \) is reached, we access its closest ancestor, which also has an outgoing edge with label \( S_i \), and splice its rectangle. As such descendants of a node are visited from left to right, we always know which part of the rectangle will be spliced next, hence each such splice takes \( O(1) \) time leading to computing \( O(n) \) such modified rectangles in \( O(n) \) total time.

In order to finalize the construction of our data structure, we compute the set of modified rectangles of one of the two types described above, and construct for them the 2D rectangle stabbing data structures for reporting (Lemma 16) and counting (Lemma 17).

**Querying.** To answer a \( \text{Report}'(j, \ell) \) or a \( \text{Count}'(j, \ell) \) query, we simply ask the corresponding 2D rectangle stabbing data structure for the point \( (L(v), \ell) = (R(v), \ell) \), where \( v \) is the node reached in \( ST_R \) from the root by conceptually reading \( S_j S_i \). In case of a reporting query, the data structure returns a set of rectangles \([x, y] \times [\ell_1, \ell_2]\) labeled with distinct values \( i \in [1, k] \). We can simply report the set of these \( i \) values. In case of a counting query, the result is simply an integer which we output. The two constructions of modified rectangles have additional nice properties however – each value \( i \) is associated with a value \( \ell_2 \). In case of the first construction, this \( \ell_2 \) is the length of the shortest suffix of \( S_i \) which is also a prefix of \( S_j \) of length at least \( \ell \); in case of the second construction, \( \ell_2 \) is the length of the longest such suffix, that is \( \ell_2 = SPL_{i,j} \).
Correctness. The correctness of the algorithm follows by the fact that point \((L(v), \ell) = (R(v), \ell)\) is enclosed by a rectangle \([L(u), R(u)] \times [0, d(u)]\) if and only if \(S_j S_i\) has a prefix of length at least \(\ell\) that is also a suffix of \(S_i\); and by the fact that the set of rectangles originating from a single \(i\) are made pairwise disjoint while their union remains unchanged.

We have thus arrived at the following lemma.

Lemma 18. For any dictionary of \(k\) strings of total length \(n\) over an integer alphabet of size \(\sigma \leq n^{O(1)}\), we can construct a data structure of \(O(n)\) words of space answering: (i) Report\((j, \ell)\) queries in \(O(\log n/\log \log n + f)\) time, where \(f\) is the size of the output; and (ii) Count\((j, \ell)\) queries in \(O(\log n/\log \log n)\) time.

By combining Lemma 18 with Observation 1 we obtain the main result of this section.

Theorem 19. For any dictionary of \(k\) strings of total length \(n\) over an integer alphabet of size \(\sigma \leq n^{O(1)}\), we can construct a data structure of \(O(n)\) words of space answering: (i) Report\((i, \ell)\) queries in \(O(\log n/\log \log n + f)\) time, where \(f\) is the output size; and (ii) Count\((i, \ell)\) queries in \(O(\log n/\log \log n)\) time.

Let us remark that the construction time for our data structures, excluding the implementation of the data structures underlying Lemmas 16 and 17, is \(O(n)\). Unfortunately, the construction time of the latter data structures (Lemmas 16 and 17) is not mentioned in [28, 45]. However, by using the construction from [15], we obtain \(O(n \log n)\) construction time, \(O(n)\) words of space, \(O(\log n + f)\) time for reporting, and \(O(\log n)\) time for counting.

Theorem 20. For any dictionary of \(k\) strings of total length \(n\) over an integer alphabet of size \(\sigma \leq n^{O(1)}\), we can construct a data structure of \(O(n)\) words of space answering: (i) Report\((i, \ell)\) queries in \(O(\log n + f)\) time, where \(f\) is the output size; and (ii) Count\((i, \ell)\) queries in \(O(\log n)\) time. The data structure construction time is \(O(n \log n)\).

Let us also remark that Report\((i, 0)\) (with the second construction of disjoint rectangles) actually answers any One-to-All\((i)\) query within the same asymptotic time: \(O(\log n + f) = O(\log n + k) = O(k)\) as \(k \geq \sqrt{n}\). While the data structure for answering Report queries occupies \(O(n)\) words of space, like the data structure for One-to-All queries, the construction time for the former is more expensive – and it is likely much slower in practice.

6 Answering Top Queries

Recall that a Top\((i, K)\) query returns exactly \(K\) elements \(j\) for which SPL\(_{i,j}\) is the largest, breaking ties arbitrarily. In case we are given an additional bound \(K' \leq k\) such that \(K \leq K'\) (e.g., we are only interested in finding \(O(1)\) many such top elements), the obvious data structure would be to store, for each \(i \in [1, k]\), the sorted list of size \(K'\) of the best answers. Such a data structure allows answering Top\((i, K)\) queries, for \(K \leq K'\), in the optimal \(O(K)\) time, but it requires \(O(kK')\) space, which for small \(K'\) may be \(O(n)\), but in general (i.e., when \(K' = k\)) leads back to the \(O(n + k^2)\)-time APSP algorithm. We show how to use our results from Section 5 to answer Top\((i, K)\) queries using \(O(n)\) space without this \(K'\) bound.

Clearly, we can assume that \(K < k\). We start by making the following crucial observation.

Observation 21. For any Top\((i, K)\) query, with \(K < k\), there exists an integer \(\ell \in [0, n - 1]\) such that Count\((i, \ell + 1)\) \(\leq K < \text{Count}(i, \ell)\).

Using the results from Section 5, we can find such an \(\ell\) in \(O(\log^2 n / \log \log n)\) time using binary search on \(\ell \in [0, n - 1]\) and the data structure for Count queries. Next we can simply compute Report\((i, \ell + 1)\) to be left with only choosing the remaining \((K - \text{Count}(i, \ell + 1))\)
elements out of all \( j \in [1, k] \) such that \( \text{SPL}_{i,j} = \ell \). Unfortunately, there can be many such elements (even \( k \)), and we do not want this to influence the query time. We have to report the remaining elements out of the ones such that \( \text{SPL}_{i,j} = \ell \) without computing or explicitly accessing all of them. Recall that, in \( ST_R \), a list of elements \( i \) such that \( S_i \) has a suffix of length exactly \( \ell \) which is also a prefix of \( S_j \) can be accessed in \( O(1) \) time after \( O(n) \)-time preprocessing by finding the ancestor of the node reached by conceptually reading \( S_j S_j \) at string depth \( \ell \) (using a WA query) and reading the first letters of its outgoing edges from left to right; since \( S_1 < \cdots < S_k \) are smaller than any element of \( \Sigma \) those values form a sorted list. Analogously, to access the list of elements \( j \) such that \( S_i \) has a suffix of length exactly \( \ell \) which is also a prefix of \( S_j \), we simply use the symmetric data structure by Observation 1.

Unfortunately, this list may contain elements \( j \) such that \( \text{SPL}_{i,j} > \ell \), and we do not want to report them again. This, however, can be fixed by maintaining a bitvector of size \( k \) as an integral part of our data structure; for each element \( j \in \text{Report}(i, \ell + 1) \), we set the \( j \)-th element of the bitvector to 1 in \( O(\text{Count}(i, \ell + 1)) = O(K) \) time. When accessing the elements of the sorted list one-by-one, we simply check if the element was already outputted using the bitvector in \( O(1) \) time. In total, we can check up to \( K \) such elements, hence the total time of merging those two parts of the output is \( O(K) \) (including the bitvector reset). We summarize the solution in Theorem 22, which is the main result of this section.

\begin{theorem}
For any dictionary of \( k \) strings of total length \( n \) over an integer alphabet of size \( \sigma \leq n^{O(1)} \), we can construct a data structure of \( O(n) \) words of space answering \( \text{Top}(i, K) \) queries in \( O(\log^2 n / \log \log n + K) \) time.
\end{theorem}

\textbf{Proof.} We start the construction of the data structure by constructing the data structures for \( \text{Report}(i, \ell) \) and \( \text{Count}(i, \ell) \) using Theorem 19. We also construct a data structure to find the list of elements \( j \) such that \( S_i \) has a suffix-prefix match of length \( \ell \) with \( S_j \) in \( O(1) \) time using Lemmas 2 and 7 and Observation 1. Finally, we also maintain a bitvector of size \( k = O(n) \). The space required by our data structure is \( O(n) \) words.

Consider a \( \text{Top}(i, K) \) query. We ask \( O(\log n) \) \( \text{Count} \) queries and a single \( \text{Report} \) query in \( O(\log^2 n / \log \log n + K) \) total time, as the output is bounded by \( K \). We index the \( \text{Report} \) result in the bitvector. We find the list (without reading its content) of elements \( j \) such that \( S_i \) has a suffix of length exactly \( \ell \) which is also a prefix of \( S_j \) in \( O(1) \) time. Finally, we access and check at most \( K \) elements from the list in \( O(K) \) total time.

The correctness of the algorithm follows by Observation 21 and Theorem 19. \hfill \blacktriangleleft

Similar to Section 5, the construction time for our data structure, excluding the implementation of Theorem 19, is \( O(n) \). If instead of Theorem 19, we employ Theorem 20, we obtain \( O(n \log n) \) construction time, \( O(n) \) words of space, and \( O(\log^2 n + K) \) query time.

\begin{theorem}
For any dictionary of \( k \) strings of total length \( n \) over an integer alphabet of size \( \sigma \leq n^{O(1)} \), we can construct a data structure of \( O(n) \) words of space answering \( \text{Top}(i, K) \) queries in \( (\log^2 n + K) \) time. The data structure construction time is \( O(n \log n) \).
\end{theorem}

References


Suffix-Prefix Queries on a Dictionary


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Merging Sorted Lists of Similar Strings

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Abstract

Merging $T$ sorted, non-redundant lists containing $M$ elements into a single sorted, non-redundant result of size $N \geq M/T$ is a classic problem typically solved practically in $O(M \log T)$ time with a priority-queue data structure the most basic of which is the simple heap. We revisit this problem in the situation where the list elements are strings and the lists contain many identical or nearly identical elements. By keeping simple auxiliary information with each heap node, we devise an $O(M \log T + S)$ worst-case method that performs no more character comparisons than the sum of the lengths of all the strings $S$, and another $O(M \log(T/\bar{e}) + S)$ method that becomes progressively more efficient as a function of the fraction of equal elements $\bar{e} = M/N$ between input lists, reaching linear time when the lists are all identical. The methods perform favorably in practice versus an alternate formulation based on a trie.

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Supplementary Material Software: https://github.com/thegenemyers/STRING.HEAP

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1 Introduction & Summary

Producing a sorted list, possibly with duplicate elements removed, from a collection of $T$ sorted input lists is a classic problem [4]. Moreover, with today’s massive data sets, where an in-memory sort would require an excessively large memory, this problem gains in importance as a component of an external, disk-based sort. Our motivating example is modern DNA sequencing projects that involve anywhere from 100 billion to 5 trillion DNA bases of data in the form of sequencing reads that are conceptually strings over the 4-letter alphabet A, C, G, T [8]. In particular, the problem of producing a sorted table of all the $k$-mers (substrings of length exactly $k$) and their counts has been the focus of much study and is used in many analysis methods for these data sets [5, 6, 7].

Priority queue implementations such as a heap, take $O(\log T)$ to extract the next minimum and insert its replacement, giving an $O(M \log T)$ merge time where $M$ is the sum of the lengths of the input lists [2]. However when the domain of the merge is strings, as opposed to say integers, then one must consider the time taken for each of the $O(\log T)$ string comparisons, which is not $O(1)$ but conceptually the average length of the longest common prefix ($lcp$) between all the compared strings. For example, this is $O(\log_{\Sigma} M)$ in the “Uniform Scenario”
Merging String Lists

where the characters of the strings are chosen with equal probability over an alphabet of size $\Sigma$. But it can be much worse, for example, when merging lists of say 21-mers each obtained from a portion of a 40X coverage DNA sequencing data set, where many strings are identical.

In the worst case, one can only assert that the time to merge the list of strings is $O(S \log T)$ where $S$ is the total number of characters in the input lists, e.g. $Mk$ for lists of $k$-mers. Assuming the Uniform Scenario, one can more accurately characterize the efficiency as $O(M \log T \log M)$ expected time. In this paper we present a method that is guaranteed to take $O(M \log T + S)$ time by modifying the heap data structure so that the amortized time spent on comparing the characters of any string while it is in the heap is never more than its length. Moreover, in the Uniform Scenario, the efficiency is $O(M(\log T + \log M))$ expected time. We call such a modified heap a string heap. Interestingly, a binary search tree augmented by a generalized list structure that also leverages lcp’s was developed by Amir et al. [1] and also achieves the bounds above, albeit with a different logic/design.

It is further true in the case of DNA sequencing data sets, that often the number of elements $N$ in the merged list is much smaller than $M$ when duplicate elements are removed. Specifically, $N$ can be as small as $M/T$ assuming the input lists themselves do not contain equal elements. With another modification to a heap, not specific to strings per se, we will achieve here an algorithm that takes $O(M \log(T/\bar{e}))$ time where $\bar{e} = M/N$ is the average number of distinct input lists a given element is in. So when all the input elements are unique the time is as usual $O(M \log T)$ but as $\bar{e}$ increases less time is taken, reaching $O(M)$ when all the input lists are identical, that is, $\bar{e} = T$. We call such a modified heap a collision heap. We show it can easily be combined with a string heap to give an $O(M \log(T/\bar{e}) + S)$ algorithm for string merging.

While the focus of this paper is on modifying a heap to support string elements, an orthogonal approach to realizing a priority queue (PQ) of strings appeared in a comprehensive paper by Thorup ([9]) that is primarily focused on integer PQs, but which in Section 6 uses a trie [3] to merge strings of, potentially large, integers in $O(M \log T + S)$ time. In bioinformatics, strings are generally over alphabets of small size $\Sigma$, e.g. 4 for DNA, so taking Thorup’s algorithm, but replacing the general integer priority queue with van Emde Boas small integer PQs [10] over domain $\Sigma$, one obtains an $O(M \log \Sigma + S)$ time algorithm. The implementation of either of these methods encounters rather larger overheads compared to simply realizing the basic approach of Thorup’s algorithm with a compact trie with $\Sigma$-element arrays for the out-edges. Moreover, because adding to the trie then becomes linear, the complexity of this simplified approach is $O(N \Sigma + S)$. Given limited values of $\Sigma$, e.g. say up to 20 for protein sequences, the trie approach is very competitive, especially for the cases where $N$ is significantly smaller than $M$.

We implemented programs to merge files of sorted strings using a regular heap, a string heap, a collision heap, a combination of the string and collision heap, and a simple compact trie and performed timing experiments on both simulated and real DNA sequencing $k$-mer data to determine their relative performance. The codes are available at github.com/thegenemyers/HEAPS. Amongst the heap-based algorithms, the string heap proves superior as the average lcp between consecutive output strings increases, and the collision heap proves superior as the collision ratio $\bar{e}$ increases. Also, the combination heap tracked the behavior of whichever of the string or collision heap proves superior, but at an overhead of roughly 5%. Against the trie approach, the string heap is faster in the uniform scenario until $T$ becomes quite large, e.g. 256 in our experiments. In scenarios where $N \ll M$ due to a uniform collision rate the trie proved fastest save for small values of $T$. For real data sets, where the collision rate is highly variable, the collision and combination heaps gave the best times. In short, the new heap methods are of both theoretical and practical interest.
## Preliminaries: Definitions and a Short Recap of Heaps

Consider \( T \) sorted lists of strings \( S_t = s_{t1}, s_{t2}, \ldots, s_{tN_t} \) of lengths \( N_t \). We assume that the elements are distinct, i.e. \( s_{tj} < s_{tj+1} \), and let \( s_{tj} = a_{jt0}a_{jt1}\ldots a_{jtn_j-1} \). Note carefully, that the first character of a string is at index 0. The problem is to produce a single sorted list \( R = r_1, r_2, \ldots, r_N \) of length \( N \) with any duplicates between the lists removed. That is, while each input list has unique strings, the same string can occur in up to \( T \) different lists. Let \( e_i \in [1, T] \) be the number of different queues the string \( r_i \) occurs in. Letting \( M = \sum_{t=1}^{T} N_t \) be the sum of the lengths of the input lists, note that \( N \) is in the range \( [M/T, M] \).

A \( T \) element heap is a complete binary tree of \( T \) nodes containing or referring to domain values to be prioritized. A heap further has the heap property when for every node, the domain values of its children are not less than its domain value. A heap can be very simply implemented as an array \( H[1..T] \) where \( H[i] \) is the datum for node \( i \), its left child is \( 2i \), and its right child is \( 2i+1 \) (if they exist, i.e. are \( \leq T \)).

In the case of merging \( T \) input lists, we will let each heap node contain the index \( t \in [1, T] \) of an input list and another array, \( V[0..T] \) will contain the current value for that list in \( V[t] \) (the role of \( V[0] \) is discussed in the next paragraph). If all the nodes greater than \( i \) have the heap property, then recall that the simple routine \( \text{Heapify}(i,x,t) \) in Figure 1 below will add the value \( x \) from list \( t \) to the heap guaranteeing that \( H \) has the heap property for all nodes greater than \( i-1 \). The routine takes time proportional to at most the height of \( i \) in the heap which is \( O(\log T) \) for all \( i \).

Let \( S[t] \) denote the \( t \)th sorted input list and assume it operates as a one-sided queue where one can \( \text{Pop} \) the next element from the list and ask if the queue is \( \text{Empty} \). We will also assume that \( \text{INFINITY} \) is an infinitely large string value greater than all those encountered as input and when a list is exhausted this place value at \( V[0] \) so that \( H[1] \) becomes 0 when all the lists are exhausted. Finally, for simplicity we assume each list has at least one element, i.e. \( N_t > 0 \) for all \( t \). Then a complete pseudo-code for the basic priority queue approach to merging \( T \) sorted lists while removing duplicate values is shown at right in Figure 1.

```c
int T
domain List S[1..T]
domain V[0..T]

Heapify(i,x,t)
{ c = i
  while (u=2c) <= T do
    if u < T and V[H[u+1]] < V[H[u]] then
      u = u+1
    if x <= V[H[u]] then break
    H[c] = H[u]
    c = u
  
  (H[c],V[t]) = (t,x)
}
```

Figure 1 The \( \text{Heapify} \) routine (left) and the overall merge algorithm (right).

In lines 1 and 2, the first element of each list is \( \text{Pop} \)’d and placed in the heap in reverse order of the nodes so that the entire heap has the heap property upon completion. The total time taken for this setup is \( O(T) \) as the sum of the heights of the nodes in a complete binary tree is of this order. Then in the while-loop of line 4, the list \( t \) with the next smallest element is \( H[1] \) and if this value is not zero (indicating the exhaustion of all the queues), then the element \( x = V[t] \) is processed in the loop body. If the value \( x \) is not a duplicate of
the last element output then it is output (lines 5 - 7). If list \( t \) is not empty then its next element replaces the element just output and the heap property is restored at node 1 (lines 8 & 11). Otherwise the element is replaced with the largest possible value \( \text{INFINITY} \) (line 9) in “queue” 0 so that when all \( T \) lists are exhausted the extraction of 0 as the queue index marks the end of the merge.

Given that \( \text{Heapify} \) takes \( O(\log T) \) time and an input element is processed with each iteration of the loop, the algorithm clearly operates in \( O(M \log T) \) time assuming domain comparisons are \( O(1) \). As discussed in the introduction this assumption is not necessarily true when the values are strings and we address this in the next section.

### 3 The String Heap

The idea for a string heap is very simple, namely, for each node also record and keep current the length of the longest common prefix between the string at the node and the string at its parent (except for the root). Let \( \text{lcp}(u, v) \) be the longest common prefix between strings \( u \) and \( v \). Then more formally, a string heap also maintains a third array \( P[1..T] \) such that \( P[i] = \text{lcp}(V[H[i]], V[H[\lceil i/2 \rceil]]) \) for \( i > 1 \). The interesting and complex part of this extension is maintaining this property during the induction of \( \text{Heapify} \) and using it to accelerate the comparison of string values by limiting the number of character comparisons involved.

Intuitively, \( \text{Heapify}(i, x, t) \) traverses the maximal left most path starting at \( i \), all of whose elements are less than \( x \) and not more than their siblings until a node \( c^* \) is reached that is either a leaf or for which all its children are not less than \( x \). The values along this path are shifted up to the node above during each iteration until \( x \) is placed at node \( c^* \) at the last. To help argue the induction to follow, it conceptually simplifies matters to think of \( x \) as being explicitly placed at the node indexed by the variable \( c \) (i.e. \( H[c] = t \)) as the algorithm descends from node \( i \) to the final placement of \( x \). From this viewpoint, at the start of each iteration of the loop of \( \text{Heapify} \), the heap satisfies the heap property at every node in the subtree rooted at \( i \) except \( c \) where \( x \) conceptually currently resides. For the array \( P \) realizing a string heap, the loop invariant is that \( P \) is correct except possibly at nodes \( 2c \) and \( 2c+1 \) as \( x \) has just been placed at their parent node \( c \). Our goal is to maintain this invariant through the next iteration of the loop where either \( x \) is found to be not greater than the children of \( c \) and the loop exits, or the algorithm descends to one of the children of \( c \) swapping \( x \) with the child’s value.

To facilitate a simpler logic around the comparison of strings, we will assume that every string ends with a special terminating character \( \$ \) that is less than any ordinary character (e.g. 0 for C-strings). With this convention, finding the \( \text{lcp} \) of two strings \( x \) and \( y \) is simply a matter of finding the first index \( \rho \) for which the strings have unequal characters or both are \$. Moreover note that \( x < y \) iff \( x[\rho] < y[\rho] \).

For all but the first iteration of the loop of \( \text{Heapify} \), note that the value that was at the current node \( c \) is now at \( \lceil c/2 \rceil \) having been exchanged with \( x \) as it has a smaller value. In what follows, we will let \( o = V[H[\lceil c/2 \rceil]] < x \) be this value and also let \( v_l = V[H[2c]] \) and \( v_r = V[H[2c+1]] \) be the strings currently at the left and right children of \( c \). Observe that it must be that \( P[2c] = \text{lcp}(o, v_l) \) and \( P[2c+1] = \text{lcp}(o, v_r) \) as these values are unchanged since the previous iteration when \( o \) was at node \( c \). Let \( p_l \) and \( p_r \) denote these values, respectively, and further let \( p = P[c] = \text{lcp}(o, x) \) in the proof/analysis that follows.

► **Theorem 1.** Given that insertions are monotone, i.e. the next value inserted is not less than the value just extracted, the \( \text{Heapify} \) routine of Figure 2 is correct once initialized. To start, it suffices to set, \( H[i] = 0 \) for all \( i \) and \( V[0] = $ \) and then perform Lines 1 and 2 of the merge given in Figure 1. After calling \( \text{Heapify} \), \( P[1] \) is the \( \text{lcp} \) of the the last value extracted and the value at the root.
Proof. First consider the situation when the heap has been correctly initialized and one is now inserting a new element $x$ as in Line 11 of the merge algorithm in Figure 1. So upon entry to Heapify, $c = i = 1$ and observe that it will be the case that $P[2]$ and $P[3]$ will have the value $\text{lcp}(o,H[V[2]])$ and $\text{lcp}(o,H[V[3]])$ where $o$ is the value that was just extracted from the root of the heap and which $x$ is now about to replace. So in order to get started we need to set $p$ to $\text{lcp}(o,x)$ where $o < x$ by the monotonicity condition. Given that nothing has been placed at $V[H[1]]$, it still has $o$ as its value, and so in Figure 2, Heapify starts correctly by setting $p$ to $\text{LCP2}(V[H[1]],x,0)$ before initiating its loop. Indeed one could imagine that $o$ is at the virtual father of the root 1.

(The reader should observe that if $o > x$, i.e. the context is not monotone, then getting the induction started also requires readjusting $P[2]$ and $P[3]$ downward to $p$ if they happen to be larger than $p$. In this case, we can no longer place a bound on the total number of character comparison made during the operation of the heap, but it will still operate correctly.)

In the case that the heap is being initialized, i.e. $i > 1$ in Line 2 of the merge algorithm, it suffices to let $o$ be the empty string, $\epsilon$, so that $P[i] = P[2i] = P[2i+1] = 0$. The conditions of the theorem correctly guarantee then that $\text{LCP2}(V[H[1]],x,0) = \text{LCP2}(S,x,0) = 0$.

We now proceed to analyze the numerous cases that arise to maintain the induction during the iterations of Heapify’s loop in terms of the relationships between the quantities $p$, $p_r$, and $p_t$. To further simplify matters observe that the treatment of the left and right children of $c$ is symmetric, so we only consider the left case, $p_r \leq p_t$, in the enumeration below knowing that the right case, $p_t < p_r$, is handled simply by exchanging the roles of left and right. Furthermore, we repeatedly use the logic that if $\text{lcp}(x,s) < \text{lcp}(s,y)$ then $\text{lcp}(x,y) = \text{lcp}(x,s)$ and $x < s$ if $x < y$.

Case 1: $p_r < p_t$ and $p < p_t$. By the case condition $\text{lcp}(x,o) = p < p_t = \text{lcp}(o,v_t)$ and since we know $x > o$ we can conclude that $x > v_t$ and $\text{lcp}(x,v_t) = \text{lcp}(x,o) = p$. Similarly $\text{lcp}(v_t,o) = p_r < p_t = \text{lcp}(o,v_t)$ and we know $v_r \geq o$ allowing us to conclude that $v_r > v_t$ and $\text{lcp}(v_r,v_t) = \text{lcp}(v_r,o) = p_r$. So $v_t$ is the smallest of $x$, $v_t$, and $v_r$ implying that the loop should descend to $2c$ with $v_t$ being placed at $c$. Moreover, $P[c]$ should be set to $p_t$, while $p$ and $P[2c+1]$ can remain unchanged having already the correct values for the next iteration.

Case 2: $p_r \leq p_t$ and $p > p_t$. By the case condition $\text{lcp}(v_t,o) = p_t < p = \text{lcp}(o,x)$ and since we know $v_t \geq o$ we can conclude that $x < v_t$ and $\text{lcp}(v_t,x) = \text{lcp}(v_t,o)$. Similarly $\text{lcp}(v_r,o) = p_r \leq p_t < p = \text{lcp}(o,x)$ and we know $v_r \geq o$ allowing us to conclude that $x < v_r$ and $\text{lcp}(v_r,x) = \text{lcp}(v_r,o)$. So the loop can terminate with $x$ being placed at node $c$. Moreover, $P[2c]$ and $P[2c+1]$ remain unchanged having yet the correct values.

Case 3: $p_r < p_t$ and $p = p_t$. First compute $p_x = p + \text{lcp}(v_t + p,x + p)$ where $s + j$ is the suffix of string $s$ beginning at position $j$. Clearly $p_x = \text{lcp}(v_t,x)$ and if $v[p_x] < x[p_x]$ then $v_t < x$, otherwise $v_t \geq x$. We have two subcases:

Subcase 3a: $v_j[p_x] < x[p_x]$. The condition $p_r < p_t$ implies $\text{lcp}(v_r,o) < \text{lcp}(o,v_t)$ and we know $v_r \geq o$ allowing us to conclude that $v_r > v_t$ and $\text{lcp}(v_r,v_t) = \text{lcp}(v_r,o)$. Thus $v_t$ is smaller than both $x$ and $v_r$. So the loop should descend to $2c$ with $v_t$ being placed at $c$. Therefore, $P[c]$ should be set to $p_t$ and $p$ to $p_x$, while $P[2c+1]$ has the correct value.

Subcase 3b: $v_j[p_x] \geq x[p_x]$. By the case conditions we know $p_r < p$ implying $\text{lcp}(v_r,o) < \text{lcp}(o,x)$ and we know $v_r \geq o$ allowing us to conclude that $v_r > x$ and $\text{lcp}(v_r,x) = \text{lcp}(v_r,o)$. Thus $x$ is not less than both $v_t$ and $v_r$. So the loop can terminate with $x$ being placed at node $c$. While $P[2c+1]$ remains correct, $P[2c]$ needs to be updated to $p_x$. 


Case 4: \( p_r = p_l \) and \( p < p_l \). Compute \( p_x = p_l + lcp(v_l + p_l, v_r + p_l) \) which is clearly \( lcp(v_1, v_r) \). If \( v_r[p_x] < v_l[p_x] \) then \( v_r < v_l \), otherwise \( v_r \geq v_l \). WLOG let’s assume \( v_l \leq v_r \), as the case \( v_r < v_l \) is symmetric. As in cases before \( p < p_l \) and \( x > o \) allow us to surmise that \( x > v_l \) and \( lcp(x, v_l) = lcp(x, o) \). Therefore \( v_l \) is smaller than \( x \) and not larger than \( v_r \), implying that the loop should descend to \( 2c \) with \( v_l \) being placed at \( c \). Therefore, \( P[c] \) should be set to \( p_l, P[2c + 1] \) to \( p_x \), while \( p \) continues to have the correct value.

Case 5: \( p_r = p_l = p \). First compute \( p_x = p + lcp3(v_l + p, v_r + p, x + p) \) where \( lcp3 \) is the 3-way common prefix, and this by the case conditions is clearly equal to \( lcp3(v_l, v_r, x) \). Therefore, \( v_l[p_x] \geq v_l[p_x] \), \( v_r[p_x] \geq v_r[p_x] \), and \( v_r[p_x] \) in direct analogy to the subcases based on the relationships between \( p, p_l, \) and \( p_r \), so we will number these 5.1, 5.2, and so on:

Subcase 5.1: \( v_r[p_x] > v_l[p_x] \) and \( x[p_x] > v_l[p_x] \). The case conditions imply \( x > v_l \) and \( v_r > v_l \) and \( lcp(v_l, x) = lcp(v_l, v_r) = p_x \). So \( v_l \) is the smallest of \( x, v_l, \) and \( v_r \) implying that the loop should descend to \( 2c \) with \( v_l \) being placed at \( c \). So \( P[c] \) should be set to \( p_l \), while \( p \) and \( P[2c + 1] \) are now clearly \( p_x \).

Subcase 5.2: \( v_r[p_x] \geq v_l[p_x] \) and \( x[p_x] \leq v_l[p_x] \). In this subcase, clearly \( x \) is not more than both \( v_l \) and \( v_r \) and \( lcp(v_l, x) = lcp(v_r, x) = p_x \). So the loop should terminate and both \( P[2c] \) and \( P[2c + 1] \) should be updated to \( p_x \).

Subcase 5.3: \( v_r[p_x] > v_l[p_x] \) and \( x[p_x] = v_l[p_x] \). First compute \( p_y = p_x + lcp(v_l + p_x, x + p) \) which is clearly \( lcp(v_l, x) \) note that the conditions to this point imply \( lcp(v_l, v_r) = lcp(x, v_r) = p_x \).

Subcase 5.3a: \( v_l[p_y] < x[p_y] \). So \( x \) is the smaller than \( x \) and \( v_r \) implying the loop should descend to \( 2c \) with \( v_l \) being placed at \( c \). So \( P[c] \) should be set to \( p_l \) and the correct new values for \( p \) and \( P[2c + 1] \) are \( p_y \) and \( p_x \), respectively.

Subcase 5.3b: \( v_l[p_y] \geq x[p_y] \). So \( x \) is not smaller than \( v_l \) and \( v_r \) implying the loop can terminate at \( c \).

Subcase 5.4: \( v_r[p_x] = v_l[p_x] \) and \( x[p_x] > v_l[p_x] \). Compute \( p_y = p_x + lcp(v_l + p_x, v_r + p_x) \) which is clearly \( lcp(v_l, v_r) \). If \( v_r[p_y] < v_l[p_y] \) then \( v_r < v_l \), otherwise \( v_r \geq v_l \). WLOG let’s assume \( v_r \leq v_r \), as the case \( v_r < v_l \) is symmetric. By the case conditions \( x > v_l \) and \( lcp(x, v_l) = p_x \). Therefore \( v_l \) is smaller than \( x \) and not larger than \( v_r \), implying that the loop should descend to \( 2c \) with \( v_l \) being placed at \( c \). Therefore, \( P[c] \) should be set to \( p_l, P[2c + 1] \) to \( p_y \), and \( p \) to \( p_x \). \( \triangle \)

Figure 2 presents the complete algorithm for the string version of Heapify embodying the case analysis above so that the \( P \)-array values are correctly maintained. Note carefully, that the \( lcp \) information in the \( P \)-array is used both to determine the relative values of the heap elements and hence direct the path that Heapify takes to insert a new element \( x \), but further also saves time on the number of character comparisons performed by only computing new \( lcp \)’s in terms of an initial \( lcp \)-offset that is common to all of the arguments to LCP2 or LCP3. So as regards complexity, the algorithm for Heapify takes \( O(\log T) \) time plus the time spent in LCP2 or LCP3 for character comparisons. Note carefully the code assumes that we are merging sorted string lists, so the value of \( x \) is not less than the value of the previous element \( o = V[H[1]] \) on the same queue \( H[1] \), i.e. the computation is monotone. We make an amortization argument to bound the total number of character comparisons as follows:
Heapify(int i, string x, int t)
{ c = i { while true do
p = LCP2(V[H[i]],x,0)
if 1 < T then
  (hr,pr) = (H[i+1],P[i+1])
else
  pr = -1
  if pr < pl then # Case 1L
    (H[c],P[c],c) = (hl,pl,1)
  else if p > pl then # Case 2L
    (H[c],P[c],P[l],c) = (hl,pl,px,l)
  else # Case 3L
    (H[c],P[c],P[l],c) = (hl,pl,px,l+1)
else
  pr = -1
  if pr > pl then # Case 1R, 2R, 3Ra, 3Rb
    (H[c],P[c],P[l],c) = (hl,pl,px,l+1)
  else if p < pl then # Case 2
    (P[l]) = px
    break
  else
    (vl,vr) = (V[hl],V[hr])
    if x[px] > vl[px] then # Case 5.2
      (P[l],P[l+1]) = (px,px)
      break
    else if vl[px] < vr[px] then # Case 5.3Rb
      (H[c],P[c],P[l],c) = (hr,pr,py,l)
      (P[l],P[l+1]) = (py,px)
      break
    else if vr[px] < vl[px] then # Case 5.3Rb
      (H[c],P[c],P[l],c) = (hr,pr,py,l)
      (P[l],P[l+1]) = (py,px)
      break
    else # Case 5.4
      if x[px] > vl[px] then # Case 5.1L
        (H[c],P[c],P[l+1],p,c) = (hl,pl,px,px,l)
      else if x[px] < vl[px] then # Case 5.2L
        (P[l]) = P[l+1] = px
        break
      else
        (H[c],P[c],P[l+1],p,c) = (hl,pl,px,px,l)
      end
    end
end
} int LCP(string x, string y, int n)
{ While true do
  (a,b) = (x[n],y[n])
  if a != b or a == $ then
    return n
    n += 1
  end
end
} int LCP3(string x, string y, string z, int n)
{ While true do
  (a,b,c) = (x[n],y[n],z[n])
  if a != b or a != c or a == $ then
    return n
    n += 1
  end
end
} Figure 2 The Heapify algorithm for a string heap.
The merge algorithm of Figure 1 when using the string heap of Figure 2 takes $O(M \log T + X)$ time where $X = |s_1| + \sum_{i=2}^{M} lcp(s_{i-1}, s_i)$ and $R^+= s_1, s_2, \ldots, s_M$ is the sequence of $M$ strings extracted from the heap over the course of the list merge, i.e. the output list if duplicates were not removed. Thus it takes $O(M \log T + S)$ worst-case time and $O(M(\log T + \log M))$ expected time in the Uniform Scenario.

**Proof.** First, observe that every string value has an lcp-value associated with it, namely, for the string $V[H[i]]$ it is $P[i]$ and it represents the number of character comparisons "charged" to its string. Examination of the case conditions reveals that when LCP2 or LCP3 is called, all the string arguments have the same lcp-value at the time of the call. Afterwords, all but one of the arguments will have its lcp-value increased to the returned value, effectively charging the comparisons of the lcp call to those arguments (NB: for LCP3 two comparisons per lcp increment are made). The total time taken then over the course of the merge is the sum of the maximum lcp-value of every string that passes through the heap. Since the lcp-value of each string is never more than the length of the string, we have our $O(S)$ bound on the total number of character comparisons.

We can more accurately characterize the number of comparisons with the observation that the maximum lcp-value that each string reaches when it is extracted from the root of the heap is its lcp with the string value extracted just before it. To see this simply review WLOG the logic involved in a value moving from node 2 to node 1 where, in all relevant cases, $P[1]$ is assigned to $p_1 = lcp(o, v_1)$ where $o$ is the last value extracted as explained previously. Further note that the comparisons for the first element extracted equals its length as its conceptual predecessor is the empty string. So the total number of character comparisons is $|s_1| + \sum_{i=2}^{M} lcp(s_{i-1}, s_i)$ over the $M$ string in $R^+$. This expression clearly reveals that the time spent comparing strings in a string heap is a function of the consecutive similarity of the strings in the final list, and immediately proves the expected time complexity claim for the Uniform Scenario as the average lcp value is $O(\log_2 M)$ in this scenario.

4 The Collision Heap

One might think that when merging sorted string lists that themselves have no duplicates, that there would be in expectation very few duplicates between the lists. This would be correct for the Uniform Scenario. But this is not true, for instance, when the problem is to merge lists of $k$-mers generated from a shotgun data set. To wit, in a coverage $c$, say 40X, data set, every part of the underlying target sequence/genome has been sampled on average 40 times and so we expect non-erroneous $k$-mers from unique parts of the target to occur on average 40 times, and a multiple thereof if from repetitive regions. So if one were to partition the data into $T$ equal sized parts, sort the $k$-mers in each part, and then merge those lists, one quite often sees the same $k$-mer in different lists. More precisely, the chance that a given $k$-mer that occurs $c$ times in the data set is not in a given input queue is $(1 - 1/T)^c$, so we expect the $k$-mer to be in $\bar{e} = T(1 - (1 - 1/T)^c) \approx T(1 - e^{-c/T})$ of the input lists. So if $T$ is say 10, then a non-erroneous, unique $k$-mer will be found in $\bar{e} = 6.5, 8.8, 9.6, \text{ or } 9.85$ of the lists if $c = 10, 20, 30, \text{ or } 40$, respectively. It was this specific use-case, that we call the "Shotgun Scenario", that motivated the development of a collision heap.

---

1 The test $x \neq \text{last}$ is simply replaced by $P[1] < |x|$ as $P[1]$ is the lcp of $x$ and the previous extracted element per Theorem 1.
The idea behind a collision heap is also very simple, namely, for each node one also records whether the value at that node is equal to its left child and its right child with a pair of boolean flags in auxiliary (bit) arrays $L[1..T]$ and $R[1..T]$. Formally, $L[i]$ has the value of the predicate $V[H[i]] = V[H[2i]]$ and $R[i]$ has the value of the predicate $V[H[i]] = V[H[2i + 1]]$. Again the interesting and somewhat complex part of this extension is maintaining these values during the induction of Heapify and using them to accelerate the handling of duplicate entries.

**Theorem 3.** The Heapify routine of Figure 3 correctly maintains the $L$ and $R$ arrays.

**Proof.** The inductive invariant for the loop of Heapify is basically that all values are correct or will be correct once the algorithm is complete, save for $H[c]$, $L[c]$ and $R[c]$ which need to be determined depending on the relative values of $x$, conceptually at $c$, and those of its current children. Let $v_l = H[V[2c]]$ and $v_r = H[V[2c + 1]]$ be the strings currently at the left and right children of $c$. There are 9 cases depending on the relative magnitudes of $x$, $v_l$, and $v_r$, where the three that entail the condition $v_l > v_r$ are treated by symmetry:

**Case 1:** $v_l < v_r$ and $x > v_l$. By the conditions, $v_l$ will move to node $c$ and the path followed descends to $2c$. $v_l < v_r$ implies that $R[c]$ should be false. However, $v_l < x$ does not imply the same for the new value of $L[c]$ as $v_l$ could be equal to the element at $2(2c)$ or $2(2c) + 1$ or both and if so, then those elements are also less than $x$ implying one or the other will replace $x$ at $2c$. Therefore $L[c]$ should be true as it will be correct and remain correct after the next loop iteration. So to recapitulate, if $L[2c] \lor R[2c]$ are true then $L[c]$ should be set to true otherwise it should be false.

**Case 2:** $v_l < v_r$ and $x = v_l$. By the conditions, the loop will terminate with $x$ finally resting at node $c$. By the case conditions it is then clear that $L[c]$ is true and $R[c]$ is false.

**Case 3:** $v_l \leq v_r$ and $x < v_l$. Again a very simple case where the loop stops and clearly $L[c] = R[c] = false$.

**Case 4:** $v_l = v_r$ and $x > v_l$. In this case, $v_l$ moves up to occupy $c$ and $x$ moves down to node $2c$. Clearly $R[c]$ should then be true as $v_l = v_r$. As argued in Case 1, if $v_l$ equals either of its children then the value of $L[c]$ needs to be true as one of these children is smaller than $x$. Otherwise $L[c]$ should be false.

**Case 5:** $v_l = v_r$ and $x = v_l$. Then the loop terminates and both $L[c]$ and $R[c]$ are true. ▶

Figure 3 presents the complete algorithm for the collision version of Heapify embodying the case analysis above so that the $L$ and $R$ array values are correctly maintained when a heap update occurs. The code is further obviously $O(\log T)$.

The value of the additional $L$ and $R$ flags is that when the top element, say $x$, is about to be extracted as the current minimum in the heap, one can find all the additional elements equal to $x$ by recursively visiting the children that are marked as equal according to the relevant $L$ and $R$ flags. In Figure 3, the routine PopHeap calls the recursive routine cohort that makes a post order traversal of the subtree of the heap of all elements equal to $x$, and places the indices of these nodes in post order in an array $G$, returning how many of them there are. Thus after calling PopHeap the array $G[1..PopHeap()]$ contains the next group of equal elements. The routine clearly takes time proportional to the number of equal elements found. The interesting part is how to replace the cohort in the heap and the time taken to do so, which we treat in the following:

**Theorem 4.** The merge algorithm of Figure 3 correctly merges the lists, outputting a unique element in each iteration and takes $O(M \log(T/\bar{c}))$ worst-case time where $\bar{c} = M/N$.  

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Merging String Lists

```java
int H[1..T]
value V[1..T]
boolean L[1..T]
boolean R[1..T]

static void Heapify(int i, value x, int t)
    { V[t] = x
      c = i
      while ((1 = (2c)) <= T)
        { hl = H[l]
          vl = V[hl]
          if (l >= T)
            vr = INFINITY
          else
            { hr = H[l+1]
              vr = V[hr]
            }
        if (vr > vl)
        { if (x > vl)
            { H[c] = hl
              L[c] = L[l] or R[l]
              R[c] = false
              c = l
            }
          else if (x == vl)
            { H[c] = t
              L[c] = true
              R[c] = false
              return
            }
        }
        else
        { if (x < vl)
            break
        }
    }

int G[1..T]

int cohort(int c, int len)
    { if (R[c])
        len = cohort(2c+1,len)
      if (L[c])
        len = cohort(2c,len)
      len += 1
      G[len] = c
      return len
    }
```

Figure 3 The `Heapify` (left) and `cohort` (upper right) and top-level merge (lower right) algorithms for a collision heap. In the main algorithm `cohort(1,0)` identifies all of the equal next elements to be output in Line 5, and then Lines 6-12 carefully replace each of these with its list successor.
Proof. While the flags allow us to easily identify the next cohort of equal elements to extract from the heap, there remains the somewhat more subtle problem of replacing all of them with their list successors. Lines 6-12 of the pseudo-code for the top-level merge at the bottom right of Figure 3 details how this is done. Because the nodes in the cohort \( G \) are in post-order, calling \textit{Heapify} on each listed node in that order guarantees a proper heap after all the elements have been replaced. In terms of complexity suppose \( e \) nodes are in the cohort for a given iteration of the loop. While the time taken in Lines 6-12 is certainly \( O(e \log T) \) we can bound this more tightly by observing that the most time is taken when the \( e \) nodes form a complete binary subtree of the heap, that is, every node has the highest height possible. In this case the lowest nodes are at height \( \log T - \log e \) and the sum over all \( e \) nodes is dominated by this as the sum telescopes (e.g. as for the time analysis for establishing the heap in Lines 1 and 2). Thus the time taken is more accurately \( O(e \log(T/e)) \). Observe that when \( e = T \) the time is \( O(e) \) and when \( e = 1 \) the time taken is \( O(\log T) \).

Looking at the overall time to produce the final list \( R \) where \( r_i \) occurs in \( e_i \) of the lists, the total time is \( O(\sum_{i=1}^{N} e_i(\log T - \log e_i)) \). By the convexity of the log-function \( \sum e_i \log_2 e_i \geq N \bar{e} \log_2 \bar{e} \) where \( \bar{e} = \sum e_i / N \) is the average value of \( e_i \). It thus follows that the total time is \( O(\sum_{i} e_i \log T - N \bar{e} \log \bar{e}) = O(M \log(T/\bar{e})) \). So when \( \bar{e} = 1 \), i.e. every input element is unique, then the time is \( O(M \log T) \) as usual. But this gradually decreases as \( \bar{e} \) approaches \( T \) where upon the time is \( O(M) \).

5 The String Collision Heap

Observing that the idea of a string heap and a collision heap are independent, one can combine the ideas obtaining an \( O(M \log(T/\bar{e}) + S) \) time algorithm. Further observe that the \( L \)- and \( R \)-arrays are not necessarily needed as \( L_{t} \) is the same as the predicate \( V[H[x]][P[x]] = \$ \) where \( x = 2c \) and \( R_{t} \) is similarly \( V[H[x]][P[x]] = \$ \) where \( x = 2c + 1 \). In words, the string of a child equals the string of its parent iff the character at its' \( lcp \)-value is the end of its' string. If one has the length \( Len[t] \) of the current string from the \( t^{th} \) input list, then the test is simply, \( P[x] = Len[H[x]] \) where \( x \) is either \( 2c \) or \( 2c + 1 \).

6 A Trie-Based Priority Queue for Strings

We briefly review trie-based implementations of a string priority queue in order to explain which approach we chose to compare against the modified heap algorithms of this paper. Given a basic Fredkin trie, adding a new string is a matter of following the path from the root of the trie spelling the common prefix with the new string, until its remaining suffix diverges at some node \( x \). A new out edge labelled with the first character of the remaining suffix is added to node \( x \) and trie nodes for the suffix are linked in. Finding the minimum string in a trie is simply a matter of following the out edge with the smallest character from each node. To delete this minimum, one finds the last node along the minimum path that has out degree greater than one, and then removes the minimum out edge from this divergent node and the suffix that follows.

If the out edges of each node are realized with a van Emde Boas priority queue for which add and delete are \( O(\log \log \Sigma) \) and finding the minimum is \( O(1) \) then adding and deleting from the queue are both \( O(\log \log \Sigma + s) \) where \( s \) is the length of the string being added or deleted. Finding the minimum element is \( O(s) \). This gives the \( O(M \log \log \Sigma + S) \) bound for the entire merge. If one further realizes a compact trie, wherein all nodes with out degree 1
are collapsed into their successor so that nodes are now labeled with string fragments, the trie is guaranteed to have $O(T)$ nodes and thus the space requirement for the trie is $O(T\Sigma)$ (excluding the space for the strings themselves).

Empirically we found that for typical values of $\Sigma$ it is actually more efficient to simply realize the out edge PQ with a $\Sigma$ element array that is directly indexed with a character. In addition, one keeps the current out-degree of the node and the current minimum out-edge. With this information finding the minimum and adding a new string is just $O(s)$. Deletion however does require traversing the out-edge array at the divergent node looking for the new minimum out-edge and so is $O(\Sigma + s)$. Offsetting this is the fact that the number of strings deleted/extracted from the trie is $N$ and not $M$, so the total complexity for this simple implementation is $O(N\Sigma + S)$ and as will be seen this empirically gives very good performance for the Shotgun Scenario.

Table 1 Performance for the Uniform Scenario.

<table>
<thead>
<tr>
<th>M, $\Sigma$, $\bar{l}p$, $\bar{e}$</th>
<th>T</th>
<th>Time (in sec.)</th>
<th>Time (in sec.)</th>
<th>Time (in sec.)</th>
<th>Time (in sec.)</th>
<th>Time (in sec.)</th>
</tr>
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<tbody>
<tr>
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<td></td>
<td>Heap</td>
<td>String Heap</td>
<td>Collision Heap</td>
<td>Combo\footnote{\textsuperscript{\textdagger}}</td>
<td>Trie</td>
</tr>
<tr>
<td>10M, 4, 10.8, 1.000</td>
<td>4</td>
<td>.654</td>
<td>.506</td>
<td>.597</td>
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7 Empirical Performance

We implemented string list merging programs using a regular heap, Heap, a string heap, Sheap, a collision heap, Cheap, a string-collision heap, SCheap, and a simple, compact trie, Trie, and measured their performance on a 2019 Mac Pro with a 2.3 GHz Intel Core i9 processor, 64GB of memory, and 8TB of SSD disk. All the codes are available at GitHub at the url github.com/thegenemyers/STRING.HEAP.
In the first set of timing experiments over synthetic data, for a given setting of parameters $M$, $T$, and $\Sigma$ we generated $T$ input files, each with $M$ 20-mers where every 20-mer over a $\Sigma$ character ASCII alphabet occurs with equally likelihood, that is, the Uniform Scenario introduced in the introduction. We chose 20 as the $k$-mer size as it is greater than the $lcp$ seen in any of the experiments. For such data we expect the $lcp$ between successive elements in the output list to be on average $\log \Sigma M$ and $\bar{e}$ to be 1 given that the average $lcp$ is less than 20 for all trials considered. In Table 1, we present timings where $\Sigma$ was set to 4, $T$ was set from 4 to 256 in steps of 2x, and $M$ was set to $10^x$ for $x = 7, 8, \text{ and } 9$. In addition, for $M = 10^9$, we also generated data sets where $\Sigma$ was also set to 8 and 16 to see the dependence of the programs, especially Trie, on $\Sigma$.

**Table 2** Performance for the Shotgun Scenario.

<table>
<thead>
<tr>
<th>$T$</th>
<th>$c$</th>
<th>$lcp$</th>
<th>$\bar{e}$</th>
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<th>Time (in sec.)</th>
<th>Time (in sec.)</th>
<th>Time (in sec.)</th>
<th>Time (in sec.)</th>
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<td>Combo'</td>
<td>Trie</td>
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The timings confirm that all algorithms are linear in $M$ and that the heap algorithms are linear in $\log T$. As $M$ becomes larger or $\Sigma$ becomes smaller the average $lcp$ between consecutive strings increases and so as expected the string heap becomes progressively faster than a regular heap. The combination heap tracks the performance of the string heap but lags by about 5% for all parameter values due to the additional overhead of maintaining information about collisions, which in these experiments basically do not occur. The trie’s behavior is basically constant, edging up slightly with $T$ due to an increase in the branching layers in the prefix of the trie. Counter intuitively, the trie becomes faster with larger $\Sigma$ as this reduces the expected number of branching layers in the trie which dominates the minor cost of searching for the smallest out-edge of a single node when deleting an entry. Thus, ultimately as $T$ increases the trie becomes the fastest, at 256 for $\Sigma = 4$, 64 for $\Sigma = 8$, and 32 for $\Sigma = 16$.

In the second set of timing experiments, we produced synthetic data sets of $k$-mers where they followed the Shotgun Scenario. We fixed $M$ at $10^9$, $k$ at 20, and then for each of $T = 4$, 8, and 16, we varied the coverage $c$ such that $c/T = .25, .5, 1.0, 1.5, 2.0, \text{ and } 3.0$. 

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As the number of collisions increases, the collision heap overtakes the string heap, with again the combination heap tracking the better of the two with an overhead of 5% or so. But our trie implementation does become faster for larger values of $T$ due to its $O(N\bar{c}\log(T/\bar{e}))$ complexity. The collision heap, in terms of $N$, has complexity $O(N\bar{c}\log(T/\bar{e}))$ which explains the behavior. Basically, the number of elements in the trie decreases rapidly from $T$ toward 1 as collisions occur greatly accelerating its operation. Nonetheless, the table reveals that for smaller values of $T$ the heap algorithms are generally superior.

The final set of experiments were for $k$-mers from a real shotgun sequencing data set, the motivating example for this work. For high-accuracy read data sets $k$, is typically chosen at 40 or more, as $k$-mers of that size are well conserved. The other difference with the synthetic Shotgun Scenario is that the $k$-mers occur with a complex frequency profile wherein some $k$-mers occur with frequency about $C$, but for example, 80% of the $k$-mers contain errors and occur once, others from a haplotype region occur roughly $C/2$ times, and so on. So in Table 3 below one will see that $\bar{e}$ is significantly less than for the synthetic examples, yet still substantially elevated. Interestingly in these cases the collision heap or combined heap perform best because they respond continuously to the collisions, beating out the string heap, and the $lcp$ is very near $k$, thus beating out the trie.

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<th>$\bar{c}$</th>
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In summary, the string heap performs best when the average $lcp$ value increases, and the collision ratio is low. The collision heap always performs better than the string heap when collisions become high. The combination heap tracks the better of the two combined methods, lagging by about 5%. The trie data structure is generally the best for large $T$ or pure collision scenarios, but on real high-fidelity shotgun data sets the collision and combination heaps proved superior.

References


PalFM-Index: FM-Index for Palindrome Pattern Matching

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Abstract
The palindrome pattern matching (pal-matching) is a kind of generalized pattern matching, in which two strings \(x\) and \(y\) of same length are considered to match (pal-match) if they have the same palindromic structures, i.e., for any possible \(1 \leq i < j \leq |x| = |y|\), \(x[i..j]\) is a palindrome if and only if \(y[i..j]\) is a palindrome. The pal-matching problem is the problem of searching for, in a text, the occurrences of the substrings that pal-match with a pattern. Given a text \(T\) of length \(n\) over an alphabet of size \(\sigma\), an index for pal-matching is to support, given a pattern \(P\) of length \(m\), the counting queries that compute the number \(\text{occ}\) of occurrences of \(P\) and the locating queries that compute the occurrences of \(P\). The authors in [I et al., Theor. Comput. Sci., 2013] proposed an \(O(n \lg n)\)-bit data structure to support the counting queries in \(O(m \lg \sigma)\) time and the locating queries in \(O(m \lg \sigma + \text{occ})\) time. In this paper, we propose an FM-index type index for the pal-matching problem, which we call the PalFM-index, that occupies \(2n \lg \min(\sigma, \lg n) + 2n + o(n)\) bits of space and supports the counting queries in \(O(m)\) time. The PalFM-indexes can support the locating queries in \(O(m + \Delta \text{occ})\) time by adding \(n \Delta \lg n + n + o(n)\) bits of space, where \(\Delta\) is a parameter chosen from \(\{1, 2, \ldots, n\}\) in the preprocessing phase.

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Keywords and phrases Palindrome matching, Generalized string pattern matching, Indexing

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1 Introduction

A palindrome is a string that can be read same backward as forward. Palindromic structures in a string are one of the most fundamental structures in the string and have been extensively studied. For example, it is known that any string \(w\) contains at most \(|w| + 1\) distinct palindromic substrings [6], and the strings reaching the maximum values have some intriguing properties [15, 28]. Another concept regarding palindromic structures is the palindrome complexity [1, 4, 2], which is the number of distinct palindromic substrings of a given length in a string.

Instead of thinking about distinct palindromic substrings, one might be interested in occurrences of palindromic substrings. The palindromic structures in such a sense are captured by the maximal palindromes from all possible “centers” in a string. Manacher’s algorithm [26], originally proposed for computing a prefix-palindrome, can be extended to compute all the maximal palindromes in \(O(|w|)\) time for a string \(w\). The authors in [18] considered the problem of inferring strings from a given set of maximal palindromes and showed that the problem can be solved in \(O(|w|)\) time.
In [19], a new concept called *palindrome pattern matching* was introduced as a generalized pattern matching. Two strings \(x\) and \(y\) of the same length are said to *palindrome pattern match* (pal-match in short) iff they have the same palindromic structures, i.e., the following condition holds: for any possible \(1 \leq i < j \leq |x| = |y|\), \(x[i..j]\) is a palindrome iff \(y[i..j]\) is a palindrome. We remark that \(x\) and \(y\) themselves are not necessarily palindromes. The palindrome pattern matching has potential applications to genomic analysis, in which some palindromic structures play an important role to estimate RNA secondary structures [21].

The pal-match problem is to search for, in a text, the occurrences of the substrings that pal-match with a pattern. Given a text \(T\) of length \(n\) and a pattern \(P\) of length \(m\), a Morris-Pratt type algorithm for solving the pal-match problem in \(O(n)\) time was proposed in [19]. The method in [19] is based on the \(l_{pal}\)-encoding of a string \(w\), denoted as \(l_{pal}w\), that is the integer array of length \(|w|\) such that \(l_{pal}w[i]\) is the length of the longest suffix palindrome of \(w[1..i]\). The \(l_{pal}\)-encoding is helpful because two strings \(x\) and \(y\) pal-match iff \(l_{pal}x = l_{pal}y\). When \(T\) is large and static, and patterns come online later, one might think of preprocessing \(T\) to construct an index for pal-matching. An index for pal-matching is to support the counting queries that compute the number \(occ\) of occurrences of \(P\) and the locating queries that compute the occurrences of \(P\). For this purpose, I et al. [19] proposed the *palindrome suffix tree* of \(T\), which is a compacted tree of the \(l_{pal}\)-encoded suffixes of \(T\). The palindrome suffix tree takes \(O(n \log n)\) bits of space and supports the counting queries in \(O(m \log \sigma)\) time and the locating queries in \(O(m \log \sigma + occ)\) time, where \(\sigma\) is the size of the alphabet from which characters in \(T\) are taken and \(occ\) is the number of occurrences.

In this paper, we present a new index, named the *PalFM-index*, by applying the technique of the FM-index [7] to the pal-match problem. In so doing we introduce a new encoding, named the \(ssp\)-encoding, that is based on the non-trivial shortest suffix-palindrome of each prefix. In contrast to the \(l_{pal}\)-encoding, the \(ssp\)-encoding has a good property to design the PalFM-index. The PalFM-index occupies \(2n \log \min(\sigma, \log n) + 2n + o(n)\) bits of space and supports the counting queries in \(O(m + \Delta occ)\) time by adding \(\Delta \log n + n + o(n)\) bits of space, where \(\Delta\) is a parameter chosen from \(\{1, 2, \ldots, n\}\) in the preprocessing phase.

### 1.1 Related work

One of the well-studied algorithmic problems related to palindromes is factorizing a string into non-empty palindromes, or in other words, recognizing a string that is obtained by concatenating a certain number of non-empty palindromes [26, 24, 12, 9, 20, 25, 3, 29]. The combinatorial properties discovered during tackling this factorization problem are useful to work on palindromes-related problems.

Developing techniques of designing space-efficient indexes for generalized pattern matching is of great interest. Our PalFM-index was inspired by that of Kim and Cho [23], which is a simplified version of the FM-index for parameterized pattern matching [13]. Indexes based on the FM-index for other generalized pattern matching problems were considered in [14, 11, 22].

### 2 Preliminaries

#### 2.1 Notations

An integer interval \(\{i, i + 1, \ldots, j\}\) is denoted by \([i..j]\), where \([i..j]\) represents the empty interval if \(i > j\).
Let $\Sigma$ be a finite alphabet, a set of characters. An element of $\Sigma^*$ is called a string. The length of a string $w$ is denoted by $|w|$. The empty string $\varepsilon$ is a string of length 0, that is, $|\varepsilon| = 0$. The concatenated string of two strings $x$ and $y$ are denoted as $x \cdot y$ or simply $xy$. The $i$-th character of a string $w$ is denoted by $w[i]$ for $1 \leq i \leq |w|$, and the substring of a string $w$ that begins at position $i$ and ends at position $j$ is denoted by $w[i..j]$ for $1 \leq i \leq j \leq |w|$, i.e. $w[i..j] = w[i]w[i+1]...w[j]$. For convenience, let $w[i..j] = \varepsilon$ if $i > j$. A substring of the form $w[1..j]$ (resp. $w[i..|w|]$) is called a prefix (resp. suffix) of $w$ and denoted as $w[..j]$ (resp. $w[..i]$) in shorthand. Note that $\varepsilon$ is a substring/prefix/suffix of any string $w$. A substring of $w$ is called proper if it is not $w$ itself. When needed we use parentheses to indicate positions in a concatenated string, for example, $(xy)[i]$ refers to the $i$-th character of the string $xy$. Hence, $(xy)[i]$ should be distinguished from $xy[i]$, which can be interpreted as the concatenated string of $x$ and $y[i]$.

Let $\prec$ denote the total order over an alphabet we consider. In particular, we will consider strings over a set consisting of integers and $\infty$, in which natural total order based on their values is employed. We extend $\prec$ to denote the lexicographic order of strings over the alphabet. For any strings $x$ and $y$ that do not match, we say that $x$ is lexicographically smaller than $y$ and denote it by $x \prec y$ iff $x[i+1] < y[i+1]$ for largest integer $i$ with $x[i..j] = y[i..j]$, where we assume that $x[i+1]$ or $y[i+1]$ refers to the lexicographically smallest character $\$ if it points out of bounds.

For any string $w$, let $w^R$ denote the reversed string of $w$, that is, $w^R = w[|w|] \cdots w[2]w[1]$. A string $w$ is called a palindromic if $w = w^R$. The radius of a palindromic $w$ is $|w|/2$. The center of a palindromic substring $w[i..j]$ of a string $w$ is $i+j\over 2$. A palindromic substring $w[i..j]$ is called the maximal palindrome at the center $i+j\over 2$ if no other palindromes at the center $i+j\over 2$ have a larger radius than $w[i..j]$, i.e., if $w[i-1] \neq w[j+1]$, $i = 1$, or $j = |w|$.

Two strings $x$ and $y$ of same length are said to palindrome pattern match (pal-match in short) iff they have the same palindromic structures, i.e., the following condition holds: for any possible $1 \leq i < j \leq |x| = |y|$, $x[i..j]$ is a palindromic iff $y[i..j]$ is a palindrome. For example, $abcbaaca$ and $bcacbbdb$ palindrome since their palindromic structures coincide (see Figure 1). Note that pal-matching induces a substring consistent equivalent relation [27], i.e., if $x$ and $y$ palindrome then $x[i..j]$ and $y[i..j]$ palindrome for any possible $1 \leq i < j \leq |x| = |y|$.

The pal-matching problem is to search for, in a text string $T$, the occurrences of the substrings that palindrome with a pattern $P$. In the pal-matching problem, an occurrence of $P$ refers to a position $i$ such that $T[i..i + |P| - 1]$ and $P$ palindrome. Throughout this paper we consider indexing a text $T$ of length $n$ over an alphabet $\Sigma$ of size $\sigma$.

**Figure 1** Illustration of the palindromic structures for pal-matching strings $abcbaaca$ and $bcacbbdb$. Check that the radii of their maximal palindromes for all possible centers, which are illustrated by two-headed arrows, coincide.
2.2 Toolbox

As a component of our PalFM-index, we use a data structure for a string \( w \) over an integer alphabet \( U \) supporting the following queries.

- \( \text{rank}_w(i, c) \): return the number of occurrences of character \( c \in U \) in \( w[i..] \).
- \( \text{select}_w(i, c) \): return the \( i \)-th smallest position of the occurrences of character \( c \in U \) in \( w \).
- \( \text{rangeCount}_w(i, j, c, d) \): return the number of the occurrences of any character in \([c..d] \subseteq U \) in \( w[i..j] \).

The Wavelet tree [17] supports these queries in \( O(\lg |\Sigma|) \) time using \( |w|\mathcal{H}_0(w) + o(|w| \lg |U|) \) bits of space, where \( \mathcal{H}_0(w) = O(\lg |U|) \) is the 0-th order empirical entropy of \( w \). The subsequent studies [8, 16] improved the complexities, resulting in the following theorem.

**Theorem 1** ([16]). For a string \( w \) over an integer alphabet \( U \), there is a data structure in \( |w|\mathcal{H}_0(w) + o(|w|) \) bits of space that supports \( \text{rank}, \text{select} \) and \( \text{rangeCount} \) in \( O(1 + \frac{\lg |U|}{\lg \lg |U|}) \) time.

We also use a data structure for the Range Maximum Queries (RMQs) over an integer array \( V \). Given an interval \([i..j]\) over \( V \), a query \( \text{RMQ}_V(i, j) \) returns a position in \([i..j]\) that has the maximum value in \( V[i..j] \), that is, \( \text{RMQ}_V(i, j) = \arg \max_{k \in [i..j]} V[k] \). We use the following result.

**Theorem 2** ([10]). For an integer array \( V \) of length \( n \), there is a data structure with \( 2n + o(n) \) bits of space that supports the RMQs in \( O(1) \) time.

2.3 FM-index

The suffix array \( \text{SA} \) of \( T \) is the integer array of length \( n + 1 \) such that \( \text{SA}[i] \) is the starting position of the lexicographically \( i \)-th suffix of \( T \).\(^1\) We define the string \( L \) (a.k.a. the **Burrows-Wheeler Transform (BWT)** [5] of \( T \)) of length \( n + 1 \) as follows:

\[
L[i] = \begin{cases} 
8 & (\text{SA}[i] = 1), \\
T'[\text{SA}[i] - 1] & (\text{SA}[i] > 1). 
\end{cases}
\]

We define the string \( F \) of length \( n + 1 \) as \( F = T'[\text{SA}[1]]T'[\text{SA}[2]] \cdots T'[\text{SA}[n + 1]] \). The so-called **LF-mapping** \( LF \) is the function defined to map a position \( i \) to \( j \) such that \( \text{SA}[j] = \text{SA}[i] - 1 \) (with the corner case \( LF(i) = 1 \) for \( \text{SA}[i] = 1 \)). A crucial point is that LF-mapping can be efficiently implemented by rank queries on \( L \) and select queries on \( F \) with \( LF(i) = \text{select}_F(\text{rank}_L(i, L[i]), L[i]) \).\(^2\) The occurrences of pattern \( P \) in \( T \) can be answered by finding the maximal interval \([P_\bullet..P_\bullet]\) in the \( \text{SA} \) array such that \( T[\text{SA}[i]..] \) is prefixed by \( P \) iff \( i \in [P_\bullet..P_\bullet] \), and computing the SA-values in the interval. For a string \( w \) and character \( c \), the so-called **backward search** computes the maximal interval in the \( \text{SA} \) prefixed by \( cw \) from that of \( w \) using a similar mechanism of the LF-mapping (see [7] for more details).

---

\(^1\) Against convention, we include the empty string that starts with the position \( n + 1 \) to \( \text{SA} \). In particular, \( \text{SA}[1] = n + 1 \) holds as the empty string is always the smallest suffix.

\(^2\) In the plain LF-mapping, select queries on \( F \) can be implemented by a simple table that counts, for each character \( c \), the number of occurrences of characters smaller than \( c \) in \( T \); but it is not the case in our generalized LF-mapping for pal-matching.
Table 1 A comparison between \( lpal \) and \( ssp \) for \( w = abbbabb \) and \( w’ = bw = abbbabbw \). The values that change when prepending \( b \) to \( w \) are underlined.

| \( w \) | a | b | b | a | b | b |
| \( lpal_{w} \) | 1 | 1 | 2 | 3 | 5 | 3 |
| \( ssp_{w} \) | \( \infty \) | \( \infty \) | 2 | 2 | 5 | 3 |
| \( w’ \) | b | a | b | b | a | b |
| \( lpal_{w’} \) | 1 | 1 | 3 | 2 | 3 | 5 |
| \( ssp_{w’} \) | \( \infty \) | \( \infty \) | 2 | 2 | 5 | 3 |

3 Encodings for pal-matching

The pal-matching algorithms in [19] are based on the \( lpal \)-encoding of a string \( w \), denoted as \( lpal_{w} \).\( lpal_{w} \) is the integer array of length \( |w| \) such that, for any position \( 1 \leq i \leq |w| \), \( lpal_{w}[i] \) is the length of the longest suffix-palindrome of \( w[1..i] \). See Table 1 for example.

Lemma 3 (Lemma 2 in [19]). For any strings \( x \) and \( y \), \( x \) and \( y \) pal-match iff \( lpal_{x} = lpal_{y} \).

Although Lemma 3 is sufficient to design suffix-tree type indexes, it seems that the \( lpal \)-encoding is not suitable to design FM-index type indexes. For example, more than one position could change when a character is prepended (see Table 1) and this unstable property makes messes up lexicographic order of \( lpal \)-encoded suffixes, which prevents us to implement LF-mapping space efficiently.

In this paper, we introduce a new encoding suitable to design FM-index type indexes for pal-matching. Our new encoding is based on the shortest suffix-palindrome for each prefix, where the shortest suffix is chosen excluding the trivial palindromes of length \( \leq 1 \). We call the encoding the shortest suffix-palindrome encoding (the \( ssp \)-encoding in short). For any string \( w \), the \( ssp \)-encoding \( ssp_{w} \) of \( w \) is the integer array of length \( |w| \) such that, for any position \( 1 \leq i \leq |w| \), \( ssp_{w}[i] \) is the length of the non-trivial shortest suffix-palindrome of \( w[1..i] \) if such exists, and otherwise \( \infty \). See Table 1 for example.

Lemma 4. Two strings \( x \) and \( y \) pal-match iff \( ssp_{x} = ssp_{y} \).

Proof. Since the \( ssp \)-encoding relies only on palindromic structures, the direction from left to right is clear.

In what follows, we focus on the opposite direction; \( x \) and \( y \) pal-match if \( ssp_{x} = ssp_{y} \). Assume for contrary that \( x \) and \( y \) does not pal-match. Without loss of generality, we can assume that there are positions \( i \) and \( j \) such that \( x[i..j] \) is a palindrome but \( y[i..j] \) is not, with smallest \( j \) if there are many. Note that the smallest assumption on \( j \) implies that \( y[i+1..j-1] \) is a palindrome: If \( y[i+1..j-1] \) is not a palindrome (clearly \( |y[i+1..j-1]| > 1 \) in such a case), \( j - 1 \) must be a smaller position that satisfies the above condition because \( x[i+1..j-1] \) is a palindrome. Let \( k = ssp_{x}[j] = ssp_{y}[j] \). Since \( x[i..j] \) is a palindrome, it holds that \( 1 < k \leq |x[i..j]| \). Moreover, \( k \neq |y[i..j]| \) as \( y[i..j] \) is not a palindrome.

Since the palindrome \( x[i..j] \) has a suffix-palindrome of length \( k \), the prefix \( x[i..i+k-1] \) of length \( k \) is a palindrome, too. On the other hand, since \( y[i..j] \) is not a palindrome that has a suffix-palindrome of length \( k \), the prefix \( y[i..i+k-1] \) of length \( k \) cannot be a palindrome. This contradicts the smallest assumption on \( j \) because \( i + k - 1 \) is a smaller position such that \( x[i..i+k-1] \) and \( y[i..i+k-1] \) disagree on their palindromic structures.

In contrast to the \( lpal \)-encoding, the \( ssp \)-encoding has a stable property when prepending a character.
Lemma 5. For any string $w$ and character $c$, there is at most one position $i$ ($1 \leq i \leq |w|$) such that $ssp_w[i] \neq ssp_w[i+1]$. Moreover, if such a position $i$ exists, $ssp_w[i] = \infty$ and $ssp_w[i+1] = i+1$.

**Proof.** By definition it is obvious that $ssp_w[i] = ssp_w[i+1]$ if $ssp_w[i] \neq \infty$. In what follows, we assume for contrary that there exist two positions $i$ and $i'$ with $1 \leq i < i' \leq |w|$ such that $ssp_w[i] = \infty > ssp_w[i+1]$ and $ssp_w[i'] = \infty > ssp_w[i'+1]$. Note that $ssp_w[i+1] = i+1$ and $ssp_w[i'+1] = i'+1$ by definition, and $(cw)[i..i+1]$ and $(cw)[i'..i'+1]$ are palindromes. Since $(cw)[i..i+1]$ is a prefix-palindrome of $(cw)[i'..i'+1]$, it is also a suffix-palindrome of $(cw)[i'..i'+1]$. It contradicts that $(cw)[i'..i'+1]$ is the non-trivial shortest suffix-palindrome of $(cw)[i..i+1]$. ▶

We consider yet another encoding based on the shortest suffix of $w[1..i-1]$ that is extended upwards when appending a character $w[i]$. The concept is closely related to the $ssp$-encoding because the extended palindrome is the non-trivial shortest suffix-palindrome of $w[1..i]$. An advantage of this new encoding is that we can reduce the number of distinct integers to be used to $O(\min(\sigma, \lg |w|))$, which will be used (in a symmetric way) to define $L_{\text{pal}}$ and obtain a space-efficient FM-index specialized for pal-matching.

For any string $w$ we partition the suffix-palindromes (including the empty suffix) by the characters they have immediately to their left and call each group a *suffix-pal-group* for $w$. We utilize the following lemma.

Lemma 6. For any string $w$, the number of suffix-pal-groups for $w$ is $O(\min(\sigma, \lg |w|))$.

**Proof.** It is obvious that the number of suffix-pal-groups is at most $\sigma$ because each character is associated to at most one suffix-pal-group. Also it is known that the lengths of the suffix-palindromes can be represented by $O(\lg |w|)$ arithmetic progressions and each arithmetic progression induces a period in the involved suffix (e.g., see [20]). Then we can see that every suffix-palindrome represented by an arithmetic progression is in the same group. Hence there are $O(\lg |w|)$ groups. ▶

The next lemma shows that pal-matching strings share the same structure of suffix-pal-groups.

Lemma 7. Let $x$ and $y$ be strings that pal-match and let $i$ and $j$ be integers with $1 \leq i < j \leq |x| = |y|$. If $x[i..i+1..]$ and $x[j+1..]$ are palindromes with $x[i] = x[j]$, then $y[i+1..]$ and $y[j+1..]$ are palindromes with $y[i] = y[j]$.

**Proof.** Since the palindrome $x[i..i+1..]$ has a suffix-palindrome of length $k = |x[j+1..]|$, it also has a prefix-palindrome of length $k$, that is, $x[i+1..i+k+1]$ is a palindrome. Also, $x[i+k+1] = x[j]$ holds. Since $x[i] = x[j] = x[i+k+1]$, $x[i..i+k+1]$ is a palindrome.

Since $x$ and $y$ pal-match, $y[i+1..]$, $y[j+1..]$ and $y[i..i+k+1]$ are palindromes. By transition of equivalence induced by the palindromes $y[i..i+k+1]$ and $y[i+1..]$, we can see that $y[i] = y[i+k+1] = y[j]$. Thus the claim holds. ▶

Let the shortest palindrome in a suffix-pal-group be the representative of the group. We assign consecutive integer identifiers starting from 1 to the suffix-pal-groups in increasing order of their representative’s lengths. See Figure 2 for example.

For any string $w$, we define the *shortest suffix-pal-group encoding* $ssp_{cw}$ of $w$ as the integer array of length $|w|$ such that, for any position $1 \leq i \leq |w|$, $ssp_{cw}[i]$ is the identifier assigned to the suffix-pal-group of the suffix-palindrome in $w[1..i-1]$ that is extended upwards by appending $w[i]$, if such exists, and otherwise $\infty$. See Table 2 and Figure 3 for example. Since
Figure 2 An example of suffix-pal-groups for babababacababacababacababa. The number enclosed in a circle denotes the pal-group-id. The suffix-palindromes in the suffix-pal-group with identifier 1 (resp. 2 and 3) have a (resp. b and c) immediately to their left. The identifiers are given in increasing order of their representative’s lengths, that is, \(|\varepsilon| = 0, |a| = 1\) and \(|ababa| = 5\).

the non-trivial shortest suffix of \(w[..i]\) is extended outwards from the representative of the suffix-pal-group for \(w[1..i − 1]\) that has \(w[i]\) immediately to the left, \(sspg_w[i]\) has essentially equivalent information to \(ssp_w[i]\). Formally the next lemma holds.

\[\textbf{Lemma 8.} \quad \text{For any string } x \text{ of length } k \text{, suppose we have the set of lengths of the representatives of suffix-pal-groups of } x[..k − 1]. \text{ Given } sspg_x[k] \text{ we can identify } ssp_x[k], \text{ and vice versa.}\]

\[\textbf{Proof.} \quad \text{It is clear that } ssp_x[k] = \infty \text{ iff } sspg_x[k] = \infty. \text{ Given } sspg_x[k] \neq \infty \text{ we can identify } ssp_x[k] \text{ from the representative of the suffix-pal-group with identifier } sspg_x[k]. \text{ Given } ssp_x[k] \neq \infty \text{ we can identify } sspg_x[k] \text{ from the representative that has length } ssp_x[k] − 2. \]

The next lemma shows that the \(sspg\)-encoding is another encoding for pal-matching, and induces the same lexicographic order with the \(ssp\)-encoding.

\[\textbf{Lemma 9.} \quad \text{Let } x \text{ and } y \text{ be strings of length } k \text{ such that } ssp_x[..k − 1] = ssp_y[..k − 1]. \text{ Then, } ssp_x[k] = ssp_y[k] \text{ iff } sspg_x[k] = sspg_y[k]. \text{ Also, } ssp_x[k] < ssp_y[k] \text{ iff } sspg_x[k] < sspg_y[k].\]

\[\textbf{Proof.} \quad \text{It follows from Lemma 7 that } x[..k − 1] \text{ and } y[..k − 1] \text{ have the same structure of suffix-pal-groups. By Lemma 8, } ssp_x[k] = ssp_y[k] \text{ if } sspg_x[k] = sspg_y[k], \text{ and vice versa. Since the identifiers of suffix-pal-groups are given in increasing order of their representative’s lengths, it holds that } ssp_x[k] < ssp_y[k] \text{ if and only if } sspg_x[k] < sspg_y[k].\]

For any string \(w\), let \(\pi(w) = sspg_{w.n}[|w|]\). Intuitively, \(\pi(w)\) holds the information from which prefix-palindrome of \(w[2..]\) the non-trivial shortest prefix-palindrome of \(w\) is extended, and the information is encoded with the identifier defined in the completely symmetric way as the case of the suffix-pal-groups. The function \(\pi(\cdot)\) will be applied to the suffixes of \(T\) to define \(F_{\text{pal}}\) and \(L_{\text{pal}}\), and the next lemma is a key to implement LF-mapping for our PalFM-index.
Table 2: A comparison between ssp<sub>w</sub> and sspg<sub>w</sub> for w = babbbabb. ssp<sub>w</sub>[6] = 5 because the non-trivial shortest suffix-palindrome of w[1..6] = babba is abba, which is of length 5. On the other hand, sspg<sub>w</sub>[6] = 2 because the shortest suffix-palindrome abba ending at 6 is extended from bbb and the suffix-pal-group to which bbb belongs for w[1..5] = babbb has the identifier 2.

\[ w = b a b b b a b b \]
\[ ssp_w = \infty \infty 3 2 2 5 3 2 \]
\[ ssdg_w = \infty \infty 2 1 1 1 2 2 \]

w = b a b b b a b b

\[ ssdg_w[6] = 2 \]

Figure 3: Illustration to show ssdg<sub>w</sub>[6] = 2 for w = babbbabb.

Lemma 10. Let x and y be strings of length \( \geq 1 \) such that \( \pi(x) = \pi(y) \). Then, ssp<sub>x</sub> ≺ ssp<sub>y</sub> iff ssp<sub>x[2..]</sub> ≺ ssp<sub>y[2..]</sub>.

Proof. Let \( i \) be the largest integer such that \( x[2..i] \) and \( y[2..i] \) pal-match. Since \( \pi(x) = \pi(y) \), using Lemma 9 in a symmetric way, it holds that \( x[2..i] \) and \( y[2..i] \) pal-match. Recall Lemma 5 that at most one \( \infty \) in ssp<sub>x[2..]</sub> (resp. ssp<sub>y[2..]</sub>) turns into the largest possible integer at the changed position when prepending \( x[1] \) (resp. \( y[1] \)). We analyze the cases focusing on the changed positions:

1. The claim clearly holds if neither ssp<sub>x</sub> nor ssp<sub>y</sub> has the changed position less than or equal to \( i + 1 \).
2. If both of ssp<sub>x</sub> and ssp<sub>y</sub> have the changed position at \( j \leq i + 1 \), it holds that ssp<sub>x</sub>[j] = ssp<sub>y</sub>[j] = \( j \) and ssp<sub>x[2..]</sub>[j - 1] = ssp<sub>y[2..]</sub>[j - 1] = \( \infty \), which also indicates that \( j < i + 1 \). Since this change does not affect the lexicographic order, the claim holds. See the left part of Figure 4 for an illustration of this case.
3. Assume ssp<sub>y</sub> has the changed position at \( j \leq i + 1 \), but ssp<sub>x</sub> does not. Since \( x[2..i] \) and \( y[2..i] \) pal-match, \( j \) cannot be less than \( i + 1 \), and hence, \( j = i + 1 \) and ssp<sub>x</sub>[i + 1] = ssp<sub>y</sub>[i + 1] < ssp<sub>x[2..]</sub>[j] < \( i + 1 \). Note that the lexicographic order between ssp<sub>x</sub> and ssp<sub>y</sub> (resp. ssp<sub>x[2..]</sub> and ssp<sub>y[2..]</sub>) is determined by that between ssp<sub>x</sub>[i + 1] and ssp<sub>y</sub>[i + 1] (resp. ssp<sub>x[2..]</sub>[i] and ssp<sub>y[2..]</sub>[i]). Since the lexicographic order between ssp<sub>x</sub>[i + 1] and ssp<sub>y</sub>[i + 1] is the same as that between ssp<sub>x[2..]</sub>[i] and ssp<sub>y[2..]</sub>[i], the claim holds. See the right part of Figure 4 for an illustration of this case.

Thus, we conclude that the lemma holds.
4 Computational results for new encodings

In this section, we show that the ssp- and ssplg-encodings can be computed in linear time for a given string.

We use the following known results.

- **Lemma 11** ([26]). For any string w, we can compute all the maximal palindromes in $O(|w|)$ time.

- **Lemma 12** (Lemma 3 in [19]). For any string w, we can compute $l_{pal}^w$ in $O(|w|)$ time.

Using Lemmas 11 and 12, we obtain:

- **Lemma 13**. For any string w, we can compute $ssp_w$ in $O(|w|)$ time.

**Proof.** Manacher’s algorithm [26] can compute the radius of the maximal palindrome in increasing order of centers in linear time. It can be extended to compute the length $l_{pal}^w[i]$ of the longest palindrome ending at each position $i$ because the maximal palindrome with the smallest center that ends at position $\geq i$ gives us the longest suffix-palindrome ending at $i$ by truncating the palindrome at $i$ (e.g., see Lemma 3 of [19]). In a similar way, we can compute the length $l_{pal}^{\prime w}[i]$ of the second longest palindrome ending at $i$.

In the third case, we use the fact that the non-trivial shortest suffix-palindrome ending at $i$ has length $\leq l_{pal}^{\prime w}[i]$ and it ends at $i - l_{pal}^w[i] + l_{pal}^{\prime w}[i]$, too.

Clearly all can be done in $O(|w|)$ time.

For any string w, let $G_w$ denote the array of length $|w|$ such that $G_w[i]$ stores the number of suffix-pal-groups for $w[1..i]$.

- **Lemma 14**. For any string w, we can compute $G_w$ in $O(|w|)$ time.
The PalFM-index of \( w \) is defined as the FM-index of \( w \) with \( \text{lpal}_w \) and \( \text{sspg}_w \) added. We only need to conceptually sort the suffixes of \( w \) in lexicographic order of their \( \text{sspg}_w \)-encoded order. We use the integer array \( \text{pal}_w \) of length \( n+1 \) such that \( \text{pal}_w[i] \) is the starting position of the \( i \)-th maximal palindrome of \( w \) and \( \text{pal}_w[1] = 1 \).

**Proof.** Let \( \text{sspg}_w \) be the array defined in a symmetric way of \( \text{ssp}_w \) such that \( \text{sspg}_w[i] \) stores the length of the non-trivial shortest prefix-palindrome starting at \( i \) (or \( \infty \) if such a palindrome does not exist). Using Lemma 13 in a symmetric way, we can compute \( \text{sspg}_w \) in \( O(|w|) \) time.

Let us focus on the palindromes involved in \( G_w[j] \). First, there is a suffix-pal-group for \( w[..j] \) that has \( w[j+1] = c \) immediately to their left. The right figure illustrates the case with \( \text{sspg}_w[i-1] \leq |w[i-1..j]| \), in which we see that the maximal palindrome \( w[i..j] \) is not the representative because there is a shorter palindrome that ends at \( j \) and has the same character \( c' \) immediately to the left.

![Figure 5](image)

**Figure 5** The left figure illustrates the case with \( \text{lpal}_w[j+1] > 1 \), in which we see that there is a suffix-pal-group for \( w[..j] \) that has \( w[j+1] = c \) immediately to their left. The right figure illustrates the case with \( \text{sspg}_w[i-1] \leq |w[i-1..j]| \), in which we see that the maximal palindrome \( w[i..j] \) is not the representative because there is a shorter palindrome that ends at \( j \) and has the same character \( c' \) immediately to the left.

**Lemma 15.** For any string \( w \), we can compute \( \text{sspg}_w \) in \( O(|w|) \) time.

**Proof.** We modify the algorithm presented in the proof of Lemma 14 slightly. Now the task is to count, for every position \( j + 1 \), the number of suffix-pal-groups for \( w[..j] \) whose representative is shorter than \( \text{ssp}_w[j + 1] - 1 \) because the number is exactly \( \text{sspg}_w[j + 1] \) by definition. We check every maximal palindrome \( w[i..j] \) and assign it to its ending position \( j \) if \( \text{sspg}_w[i] \leq |w[i-1..j]| \) and \( \text{sspg}_w[j + 1] - 1 > j - i + 1 \). Finally the number of representatives assigned to \( j \) plus one is \( \text{sspg}_w[j + 1] \). Similarly to the proof of Lemma 14, all can be done in \( O(|w|) \) time.

**5 PalFM-index**

The PalFM-index of \( T \) conceptually sort the suffixes of \( T \) in lexicographic order of their \( \text{ssp}-\text{encodings} \) (or equivalently \( \text{sspg}-\text{encodings} \)). Let \( \text{SA}_\text{pal} \) be the integer array of length \( n + 1 \) such that \( \text{SA}_\text{pal}[i] \) is the starting position of the \( i \)-th suffix of \( T \) in \( \text{ssp}-\text{encoded} \) order. We define the strings \( F_\text{pal} \) and \( \text{l}_\text{pal} \) of length \( n + 1 \) based on \( \pi \) function applied to the sorted suffixes. Formally, for any position \( i \) (\( 1 \leq i \leq n + 1 \)) we define:
Thus we focus on a single step of backward search. In a general setting, for any string $w$ with $\sigma$ distinct symbols in $\Sigma$, we show how to compute $\pi$-interval refer to the maximal interval $[b..e]$ that pal-matches with $w$ iff $i \in [b..e]$. A single step of backward search computes $cw$-interval from $w$-interval for some character $c$.

The following theorems are the main contributions of this paper.

**Theorem 16.** Let $T$ be a string of length $n$ over an alphabet of size $\sigma$. There is a data structure of $2n \log \min(\sigma, \lg n) + 2n + o(n)$ bits of space to support the counting queries for the pal-matching problem in $O(m)$ time, where $m$ is the length of a given pattern $P$.

**Proof.** We use the data structures of Theorem 1 for $L_{pal}$ and $F_{pal}$, and the RMQ data structure of Theorem 2 for the integer array $V$ with $V[i] = LF_{pal}(i)$. Since the number of distinct symbols in $L_{pal}$ and $F_{pal}$ are $O(\min(\sigma, \lg n))$ by Lemma 6, the data structures occupy $2n \log \min(\sigma, \lg n) + 2n + o(n)$ bits of space in total and all queries (rank, select, rangeCount and RMQ) can be supported in $O(1)$ time.

The number of occurrences of $P$ can be answered by computing the width of $P$-interval. Thus we focus on a single step of backward search. In a general setting, for any string $w$ and a character $c$, we show how to compute $cw$-interval $[b'..e']$ in $O(1)$ time from $w$-interval $[b..e]$, $\pi(cw)$ and the number $g$ of prefix-pal-groups of $w$. The procedure differs depending on $\pi(cw) = \infty$ or not.
1. When \( \pi(cw) = k \neq \infty \). Using Lemma 9 in a symmetric way, \([b',e']\) is obtained by mapping the positions of \( \pi(cw) \) in \( \text{l}_\text{pal}[b,e] \) by the \( \text{l}_\text{pal} \) function. More specifically, \( b' = \text{select}_{\text{l}_\text{pal}}(\text{rank}_{\text{l}_\text{pal}}(b - 1,k) + 1,k) \) and \( e' = \text{select}_{\text{l}_\text{pal}}(\text{rank}_{\text{l}_\text{pal}}(e,k),k) \), which can be computed in \( O(1) \) time.

2. When \( \pi(cw) = \infty \). We note that \([b',e']\) is the maximal interval such that \( T[\text{SA}_\text{pal}[i]..] \) does not have non-trivial prefix-palindrome (i.e. \( \pi(T[\text{SA}_\text{pal}[i]..]) = \infty \)) or \( T[\text{SA}_\text{pal}[i]..] \) has the non-trivial shortest prefix-palindrome of length longer than \( |cw| \) (i.e. \( \pi(T[\text{SA}_\text{pal}[i]..]) > g \)). Thus, \( e' - b' + 1 \) is equivalent to the number of occurrences of values larger than \( g \) in \( \text{l}_\text{pal}[b,e] \), which can be computed in \( \text{rangeCount}_{\text{l}_\text{pal}}(b,e,g,\infty) \) in \( O(1) \) time. Moreover, it holds that \( e' = \text{LF}_{\text{pal}}(\text{RMQ}_{\text{l}_\text{pal}}(b,e)) \) because \( \text{ssp}(T[\text{SA}_\text{pal}[i]..]) \) with \( \pi(T[\text{SA}_\text{pal}[i]..]) = \text{l}_\text{pal}[i] > g \) is always lexicographically larger than \( \text{ssp}(T[\text{SA}_\text{pal}[j]..]) \) with \( \pi(T[\text{SA}_\text{pal}[j]..]) = \text{l}_\text{pal}[j] \leq g \). Thus, we can compute \([b',e']\) in \( O(1) \) time.

Backward search for \( P \) requires \( \pi(P[i..]) \) and the number \( g \) of prefix-pal-groups of \( P[i..] \) for all \( 1 \leq i \leq m \), which can be computed by \( \text{sspg}_{P_i} \) and \( G_{P_i} \) in \( O(m) \) time using Lemmas 15 and 14.

Putting all together, we get the theorem.

\[\text{Theorem 17.} \quad \text{Let} \ T \ \text{be a string of length} \ n \ \text{over an alphabet of size} \ \sigma \ \text{and} \ \Delta \ \text{be an integer in} \ [1..n]. \ \text{There is a data structure of} \ 2n\lg \min(\sigma,\lg n) + \frac{n}{\Delta} \lg n + 3n + o(n) \ \text{bits of space to support the locating queries for the pal-matching problem in} \ O(m + \Delta \text{occ}) \ \text{time, where} \ m \ \text{is the length of a given pattern} \ P \ \text{and} \ \text{occ} \ \text{is the number of occurrences to report.}\]
Proof. We use the data structure and the algorithm of Theorem 16 to compute \( P \)-interval in 
\( 2n(1 + \lg \min(\sigma, \lg n)) + o(n) \) bits of space and \( O(m) \) time. The occurrences of \( P \) (in the sense of pal-matching) can be answered by the \( SA_{\text{pal}} \)-values in \( P \)-interval. We employ exactly the same sampling technique used in the FM-index to retrieve \( SA_{\text{pal}} \)-values (e.g., see [7]): We make a bit vector \( B \) of length \( n + 1 \) marking the positions \( i \) in \( SA_{\text{pal}} \) such that \( SA_{\text{pal}}[i] = \Delta k + 1 \) for some integer \( k \), and the sparse suffix array \( S \) holding only the marked \( SA_{\text{pal}} \)-values in the order. \( B \) is equipped with a data structure to support the rank queries and the additional space to Theorem 16 is 
\( \frac{\Delta}{2} \lg n + n + o(n) \) bits in total.

If position \( i \) is marked, \( SA_{\text{pal}}[i] \) is retrieved by \( S[\text{rank}_{B}(i, 1)] \) in \( O(1) \) time. If position \( i \) is not marked, we apply LF-mapping \( k \) times from \( i \) until we reach a marked position \( j \) and retrieve \( SA_{\text{pal}}[i] \) by \( S[\text{rank}_{B}(j, 1)] + k \). Since text positions are marked every \( \Delta \) positions, the number \( k \) of LF-mapping steps is at most \( \Delta \), and hence, \( SA_{\text{pal}}[i] \) can be retrieved in \( O(\Delta) \) time. Therefore we can report each occurrence of \( P \) in \( O(\Delta) \) time, and the theorem follows.

6 Conclusions and future work

In this paper, we developed new encoding schemes for pal-matching and proposed the PalFM-index, a space-efficient index for pal-matching based on the FM-index. Future work includes to present an efficient construction algorithm of the PalFM-index, and to reduce the space requirement (e.g. by incorporating with the idea of [13]). Another interesting research direction would be to develop a general framework to design FM-index type indexes in generalized pattern matching. We believe that switching encoding from \( lpal \) to \( ssp \) to design the PalFM-indexes gives a good hint to pursue this direction, and conjecture that any generalized pattern matching under a substring consistent equivalent relation [27] admits such shortest positional encodings to design FM-index type indexes.

References


Computing MEMs on Repetitive Text Collections

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Abstract

We consider the problem of computing the Maximal Exact Matches (MEMs) of a given pattern $P[1..m]$ on a large repetitive text collection $T[1..n]$, which is represented as a (hopefully much smaller) run-length context-free grammar of size $g_{rl}$. We show that the problem can be solved in time $O(m^2 \log^3 n)$, for any constant $\epsilon > 0$, on a data structure of size $O(g_{rl})$. Further, on a locally consistent grammar of size $O(\delta \log^2 n)$, the time decreases to $O(m \log m (\log m + \log^\epsilon n))$. The value $\delta$ is a function of the substring complexity of $T$ and $\Omega(\delta \log^2 n)$ is a tight lower bound on the compressibility of repetitive texts $T$, so our structure has optimal size in terms of $n$ and $\delta$.

1 Introduction and Related Work

Mutations and experimental sequencing errors make exact pattern matching seldom used in Bioinformatic applications, except possibly for very short patterns and some niche applications [19, 37, 28]. A much more interesting problem is that of finding the Maximal Exact Matches (MEMs) of a given pattern $P[1..m]$ in a text $T[1..n]$. A MEM is a maximal substring $P[i..j]$ that appears in $T$ (i.e., $P[i-1..j]$ and $P[i..j+1]$ are out of bounds or do not occur in $T$). This is useful, for example, to find long conserved areas of a gene or to best align a read (where $m$ is typically in the hundreds or thousands) on a reference genome (where $n$ can be in the billions), and even to find similarities between two genomes. In this paper we are interested in the case where $T$ is known in advance and can be indexed.

Finding MEMs is a classic problem in stringology and can be solved in optimal $O(m)$ time using a suffix tree of $T$ [41, 31] (see, e.g., the similar problem of computing matching statistics [19, Sec. 7.8]). Suffix trees, even if using linear space, are too large to maintain in main memory for current text collection sizes, however. The suffix tree of a single human genome, for example, with $n \approx 3 \cdot 10^9$, may take 60GB with a decent implementation. This makes suffix trees hard to use directly on current bioinformatic collections. Even if lower-space alternatives can replace suffix trees for most tasks [28, MEMs in Sec. 11.1.3], this space reduction is still insufficient to face current projects for sequencing millions of human genomes (see https://b1mg-project.eu).

A fortunate situation is that many of the fastest growing text collections are highly repetitive [33]. For example, collections of genomes of the same species feature a small percentage of differences between any pair of genomes. Several text indices exploiting repetitiveness to reduce space have appeared [34]. Those indices may take orders of magnitude less space than the raw data, and even more orders less space than a suffix tree on the data.

Those compressed indices support exact pattern matching, that is, they can list all the positions where $P$ occurs in $T$. While useful, this is less than the full suffix tree functionality, and insufficient to efficiently implement the classic $O(m)$-time MEM finding algorithm.
Compressed suffix trees for highly repetitive text collections do exist, but do not compress that much. Gagie et al. [17] show how to simulate a suffix tree within space $O(r \log \frac{n}{r})$, where $r$ is the number of equal-letter runs in the BWT [7] of $T$. It could find the MEMs in time $O(m \log \frac{n}{r})$ if we run the algorithm backwards on $P$, using operations parent and Weiner link instead of child and suffix link. The problem is the space: while $r$ is an accepted measure of repetitiveness [21], it is a weak one [33, 21], and multiplying it by $\log \frac{n}{r}$ makes it grow by an order of magnitude. Current implementations of compressed suffix trees for repetitive texts achieve remarkable space, but still use at least 2–4 bits per symbol [39, 15, 8, 5].

Another trend has been to expand the functionality of a more basic compressed text index for repetitive texts so as to support specific operations, MEMs in our case. Bannai et al. [1] show how to compute matching statistics (from where MEMs are easily extracted in $O(m)$ time) by extending the RLBWT-index [29], in $O(m(s + \log \log n))$ time and $O(r)$ space, with the help of a data structure that provides access to a symbol of $T$ in time $O(s)$. This can be, for example, the samples of the RLBWT-index, which add $O(n/s)$ space to the index, or a context-free grammar of $T$, which provides access in time $s = O(\log n)$ [4]. Various implementations of this idea [38, 6, 40] showed its practicality on large genome collections, with indices that are an order of magnitude smaller than the text.

All those results have been obtained on the so-called suffix-based compressed indices for repetitive collections [34]. This is natural because those emulate variants of suffix trees or arrays [30], which simplifies the problem of simulating the suffix tree traversal of the classic MEM-finding algorithm. Even the naïve algorithm of searching for all the $O(m^2)$ substrings of $P$ can be run in $O(m^2 \log \log n)$ time on those $O(r)$-sized indices.

The problem is much harder on the so-called parsing-based indices [34]. Those are potentially smaller than the suffix-based indices because they build on stronger measures of repetitiveness. For example, the size $g$ of the smallest context-free grammar that generates $T$ is usually considerably smaller than $r$ [33]. Because these indices cut $T$ into phrases, even exact pattern matching is complicated because the occurrences of $P$ can appear in many different forms, and many possible cuts of $P$ must be tried out ($m - 1$ in the general case) [12]. This makes the problem of finding MEMs considerably harder. We are only aware of the results of Gao [18], who computes matching statistics in time $O(m^2 \log^\epsilon \gamma + m \log n)$ using $O(\delta \log \frac{n}{\epsilon})$ space (for any constant $\epsilon > 0$), or $O(m^2 + m \log \gamma \log \log \gamma + m \log n)$ using $O(\delta \log \frac{n}{\epsilon} + \gamma \log \gamma)$ space. Here $\delta \leq \gamma$ are lower-bounding measures of repetitiveness [22, 11]. The size $O(\delta \log \frac{n}{\epsilon})$ matches a tight lower bound on the size of compressed representations of $T$ [25], so a structure of this size uses asymptotically optimal space for every $n$ and $\delta$.

Let $g_{rl}$ be the size of any run-length context-free grammar generating $T$ (those include and extend classic context-free grammars). The smallest such grammar is of size $g_{rl} = O(\delta \log \frac{n}{\epsilon})$ [25]. We first show that, on an index of size $O(g_{rl})$, one can compute the MEMs in time $O(m^2 \log \epsilon g_{rl})$, for any constant $\epsilon > 0$. This is done by sliding the window $P[i \ldots j]$ of the classic algorithm while we simulate the process of searching for that window with the grammar. The simulation is carefully crafted to avoid expensive operations, so the time stays proportional to the number of cuts tried out on a single search for $P$. The space $O(g_{rl})$ is the least known to support direct access to $T$ with logarithmic time guarantees [33]. The result essentially matches the first one of Gao, which could also run within $O(g_{rl})$ space.

We further show that, on a particular grammar featuring local consistency properties [24], we can reduce the time to $O(m \log m(\log m + \log^\epsilon n))$ by exploiting the fact that only $O(\log(j - i + 1))$ cuts need to be tried out for $P[i \ldots j]$, and using much more sophisticated techniques to amortize the costs. This grammar is of size $O(\delta \log \frac{n}{\epsilon})$, optimal for every $n$ and $\delta$, and within this space we sharply break the quadratic time of previous solutions.
2 Maximal Exact Matches (MEMs) and How to Find Them

We assume the usual notation on strings $S[1..n]$ and that the reader is familiar with the concepts related to suffix trees [41, 31, 13]. We start by defining MEMs.

- **Definition 1.** A Maximal Exact Match (MEM) of a pattern $P[1..m]$ in a string $T$ is a substring $P[i..j]$ that occurs in $T$, but in addition
  - $i = 1$ or $P[i-1..j]$ does not occur in $T$, and
  - $j = m$ or $P[i..j+1]$ does not occur in $T$.

- **Definition 2.** Given a text $T[1..n]$ that can be preprocessed, the MEM-finding problem is that of, given a pattern $P[1..m]$, return the range $(i,j)$ of each of its MEMs $P[i..j]$ in $T$, in increasing order of $i$ (or $j$). A position where each MEM occurs in $T$ must also be returned.

The MEM finding problem can be solved in $O(m)$ time with a suffix tree. Algorithm 1 shows how, abstracting away some complications of implementing it on the long edges of suffix trees. The next problem is strongly related to the MEM finding problem.

- **Definition 3.** Given a text $T[1..n]$ that can be preprocessed, the matching statistics problem is that of, given a pattern $P[1..m]$, return the length $M[k]$ of the longest prefix of $P[k..]$ that occurs in $T$, for every $1 \leq k \leq m$. A position where each such longest prefix occurs must be given for each $k$.

Given a solution to the MEM finding problem, $(i_1,j_1),\ldots,(i_s,j_s)$, we compute the matching statistics as follows. Set all $M[k]$ to zero and then traverse the tuples $(i_r,j_r)$ in order. Set $M[k] = j_r - k + 1$ for all $i_r \leq k \leq \min(j_r,i_{r+1}-1)$, assuming $i_{r+1} = m + 1$. The occurrence of each $M[k] > 0$ is that of its $(i_r,j_r)$ shifted by $k-i_r$. Conversely, given the matching statistics $M[k]$ for $1 \leq k \leq m$, we obtain the MEMs by reporting, for increasing $i$, every pair $(i,i + M[i] - 1)$ such that $i = 1$ or $M[i] \geq M[i-1]$, and $M[i] > 0$. Therefore, both problems are interchangeable as one can convert one output to the other in optimal $O(m)$ time. Gusfield [19, Sec. 7.8] shows how to compute matching statistics with the suffix tree.

---

**Algorithm 1** Finding the MEMs of $P[1..m]$ in $T$ using the suffix tree of $T$.

```plaintext
1 \hspace{1em} i \leftarrow 1; j \leftarrow 0;
2 \hspace{1em} v \leftarrow \text{suffix tree root};
3 \hspace{1em} \text{while } j < m \text{ do}
4 \hspace{2em} if \hspace{1em} v \text{ has no child labeled } P[j+1] \text{ then}
5 \hspace{3em} i \leftarrow i + 1; j \leftarrow j + 1;
6 \hspace{2em} \text{end}
7 \hspace{2em} \text{else}
8 \hspace{3em} \text{while } j < m \text{ and } v \text{ has a child labeled } P[j+1] \text{ do}
9 \hspace{4em} j \leftarrow j + 1; v \leftarrow \text{the child of } v \text{ by } P[j+1];
10 \hspace{3em} \text{end}
11 \hspace{2em} \text{report } (i,j) \text{ with some occurrence of } v;
12 \hspace{2em} \text{while } i \leq j < m \text{ and } v \text{ has no child labeled } P[j+1] \text{ do}
13 \hspace{3em} i \leftarrow i + 1; v \leftarrow \text{the suffix link of } v;
14 \hspace{2em} \text{end}
15 \hspace{2em} \text{end}
16 \hspace{1em} \text{end}
```
### 3 Grammar based Indices

Let $T[1..n]$ be a text. Grammar-based compression of $T$ consists in replacing it by a context-free grammar (CFG) that generates only $T$ [23]. The compression ratio is then the size of the grammar divided by the text size.

We consider a slightly more powerful type of grammar called run-length context-free grammar (RLCFG), which includes run-length rules of constant size. To simplify, we disallow rules of the form $A \rightarrow \varepsilon$, which are easily removed without increasing the grammar size.

- **Definition 4.** A Run-Length Context-Free Grammar (RLCFG) for $T$ is a context-free grammar that generates (only) $T$, having exactly one rule per nonterminal $A$. The rules are of the form $A \rightarrow B_1 \cdots B_k$ for $k > 0$ and terminals or nonterminals $B_i$ (this rule is said to be of size $k$), and of the form $A \rightarrow B^k$ for $k > 1$ and a terminal or nonterminal $B$, which is identical to $A \rightarrow B \cdots B$ with $k$ copies of $B$, but is said to be of size 2. The size of the RLCFG is the sum of the sizes of all of its rules. A Context-Free Grammar (CFG) for $T$ is a RLCFG for $T$ that does not use rules of the form $A \rightarrow B^k$.

Clearly, the size $g_{rl}$ of the smallest RLCFG for $T$ is always less than or equal to the size $g$ of the smallest CFG for $T$. Grammar-based compression (with or without run-length rules) has proved to be particularly effective on highly repetitive texts [34]. While finding the smallest grammar is NP-hard [10], heuristics like RePair obtain very good results [27].

Note that our RLCFGs have a unique parse tree, defined as follows [11, Sec. 4].

- **Definition 5.** The parse tree of a RLCFG for $T$ has a root labeled with the initial symbol. If a node is labeled $A$ and its rule is $A \rightarrow B_1 \cdots B_k$, then the node has $k$ children labeled $B_1, \ldots, B_k$ left to right. If its rule is $A \rightarrow B^k$, then the node has $k$ children labeled $B$. It follows that the $i$th left-to-right leaf of the parse tree is labeled $T[i]$.

While the parse tree has size $\Theta(n)$, a convenient representation of a RLCFG is the so-called grammar tree, which is of size $O(g_{rl})$ [11, Sec. 6].

- **Definition 6.** The grammar tree of a RLCFG is obtained by pruning its parse tree, preserving the leftmost internal node labeled $A$ for each nonterminal $A$, and converting the others to leaves. Further, for the remaining internal nodes labeled $A$ with rules $A \rightarrow B^k$ we preserve their first child only, replacing the other $k - 1$ children (which are leaves) with a single special leaf labeled $B[k-1]$. If the RLCFG size is $g_{rl}$, its grammar tree has $g_{rl} + 1$ nodes.

We will sometimes identify a nonterminal with its (only) internal node in the grammar tree. We call $\text{exp}(A)$ the string of terminals to which symbol $A$ expands, and $\text{exp}(a) = a$ for terminals $a$. The grammar tree defines a parse of $T$, as follows.

- **Definition 7.** The grammar tree, with leaves $v_1, \ldots, v_k$, induces the parse $T = \text{exp}(v_1) \cdot \text{exp}(v_2) \cdots \text{exp}(v_k)$ into phrases $\text{exp}(v_i)$.

A classic grammar-based index [12] divides the occurrences of a pattern $P[1..m]$ into primary and secondary, depending on whether they cross a phrase boundary or lie within a phrase, respectively (if $m = 1$, its occurrences ending a phrase boundary are taken as primary). It uses the fact that every occurrence has primary occurrences and that all the secondary ones can be found inside pruned leaves of nonterminals that contain other occurrences. In this paper we will be interested in the mechanism to find the primary occurrences. This is based on the parsing, but defined in a particular way to avoid reporting multiple times the primary occurrences that cross several phrase boundaries. The mechanism was extended to RLCFGs [11, Sec. 6 and App. A].
Definition 8. Let $X$ and $Y$ be multisets of strings defined as follows. For each rule $A \to B_1 \cdots B_t$, for each $1 < s \leq t$, the string $\exp(B_{s-1})^{rev}$ (i.e., $\exp(B_{s-1})$ read backwards) is inserted in $X$ and the string $\exp(B_s) \cdots \exp(B_t)$ is inserted in $Y$; we say those two are corresponding strings. Similarly, for each rule $A \to B'$, $\exp(B')^{rev}$ is inserted in $X$ and $\exp(B)^{rev}$ (i.e., $t-1$ concatenations of $\exp(B)$) is inserted in $Y$. A grid $G$ has one row per string in $Y$ and one column per string in $X$. After lexicographically sorting $X$ and $Y$, a point $(x, y)$ is set in $G$ if the $x$th string of $X$ corresponds to the $y$th string of $Y$.

The grammar-based index includes a Patricia tree $P_X$ storing the strings of $X$ and another Patricia tree $P_Y$ storing the strings of $Y$ [32]. Let us add some data to nodes for our convenience. Each Patricia tree node $v$ stores its range $[v^1, v^2]$ of the left-to-right ranks of the leaves descending from $v$. The edges of the Patricia tree nodes can represent virtual strings, so prefixes that end in the middle of an edge that leads to a node $v$ correspond to virtual nodes $u$; the range $[u^1, u^2]$ is the same $[v^1, v^2]$. The nodes $v$ also store their string depth $|v|$, which is also easily computed for virtual nodes as we descend or ascend in the Patricia tree.

Each primary occurrence consists of a suffix of some string $X \in X$ matching $P[1 \ldots i]$ corresponding to some string $Y \in Y$ whose prefix matches $P[i+1 \ldots m]$, for some $1 \leq i < m$ (if $m = 1$, it is just a suffix of $X$ matching $P$) [11, Sec. A.4]. Therefore, to find the primary occurrences of $P$, the index tries out every cutting point $i$, and searches $P_X$ for $P[1 \ldots i]^{rev}$ and $P_Y$ for $P[i+1 \ldots m]$. If both nodes $x \in P_X$ and $y \in P_Y$ exist, then the points in the orthogonal range $[x^1, x^2] \times [y^1, y^2]$ of $G$ represent the primary occurrences of $P$ cut at position $i$, and are efficiently found with a geometric data structure on $G$. By storing the position $t$ of $T$ where $\exp(B_{s-1})$ ends for such point, we know that $P$ occurs in $T[t-i+1 \ldots t-i+m]$ (the actual index stores pointers to the grammar tree, but this suffices for us).

Both the Patricia trees and the grid take $O(g_{rt})$ space. The index also needs to verify the matches of the Patricia trees. It uses an $O(g_{rt})$-space data structure $A$ that can read, in $O(\ell)$ time, any length-$\ell$ prefix or suffix of $\exp(A)$, for any nonterminal $A$ [11, Lem. 6.6].

The generic idea follows that of Algorithm 1, sliding a window $P[i \ldots j]$ along the pattern. We maintain a set of so-called active positions $r \in [i \ldots j]$.

Definition 9. A position $r \in [i \ldots j]$ is active if $P[r+1 \ldots j]$ prefixes some string in $P_Y$.

Note that, since we slide the window $P[i \ldots j]$ forwards, once a position $r$ becomes inactive, it will not become active again.

4 A Quadratic-Time Solution

We now present a quadratic-time solution that works with any RLCFG of size $g_{rt}$ for $T$; we use the $O(g_{rt})$-space data structures described in the previous section. Since any CFG is a particular case of RLCFG, our algorithm also runs with any CFG.

The generic idea follows that of Algorithm 1, sliding a window $P[i \ldots j]$ along the pattern. We maintain a set of so-called active positions $r \in [i \ldots j]$.

Algorithm

The algorithm maintains the invariant that, when the window is $P[i \ldots j]$, $(i, j)$ is the last MEM of $P[1 \ldots j]$ (if $i \leq j$) and all the MEMs ending before $j$ have already been reported. It maintains the set $R \subseteq [i \ldots j]$ of active positions, and for each such active position $r \in R$:  

4.1 Algorithm

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The node \( y_r \in P_Y \) corresponding to \( P[r + 1..j] \); this node can be virtual. Note that \([y_1^r, y_2^r] \) is the same range of rows in \( G \) of the strings of \( Y \) that start with \( P[r + 1..j] \).

The length \( \ell_r \) of the maximum prefix of \( P[r + 1..] \) that prefixes a string in \( P_Y \); note that \( r \) is active iff \( r + \ell_r \geq j \).

The node \( x_r \in P_X \) corresponding to the longest prefix of \( P[i..r]^{rev} \) that exists in \( P_X \), and such that there are points in \( G \) in the range \([x_1^r, x_2^r] \times [y_1^r, y_2^r] \). Note again that \( x_r \) can be virtual and that \([x_1^r, x_2^r] \) is the same range of columns in \( G \) of the strings of \( X \) that start with \( P[r - |x_r| + 1..r]^{rev} \).

Our algorithm, depicted in Algorithm 2, iterates over \( j \), from 0 to \( m - 1 \), and at each cycle it extends the current window to end in \( j + 1 \). When \( i = j + 1 \) (including when we start with \( i = 1 \) and \( j = 0 \)), the window is empty and there are no active positions. Line 3 first sees, in this case, if we can descend from the root of \( P_X \) by \( P[j + 1] \), to start a new nonempty substring \( P[i + 1, j + 1] \). If this is not possible, it just increases \( i \) and goes for the next value of \( j \). Otherwise, there will be active positions for the window ending at \( j + 1 \) and we enter into the main process.

Lines 5–7 first create the new active position \( r = j + 1 \), with corresponding \( y_r \) set at the root of \( P_Y \). To compute \( \ell_r \), we descend in \( P_Y \) as much as possible by \( P[r + 1..] \). To compute \( x_r \), we also descend in \( P_X \) as much as possible by \( P[i..r]^{rev} \). Those are classic Patricia tree searches, first reaching a candidate node \( v \) by comparing only the branching characters in the trie, and then verifying which ancestor of \( v \) is the correct answer. The verification proceeds by extracting the needed prefix from \( \langle v \rangle \) in \( P_Y \) (at most \( \ell_r + 1 \) characters) or the needed suffix in \( P_X \) (at most \( |x_r| + 1 \) characters).

Lines 8–16 then remove the active positions that do not reach \( j + 1 \) and updates the variables for the surviving ones. Line 10 first removes the active positions \( r \) where \( r + \ell_r = j \). On the remaining ones, each \( y_r \) moves to its child by \( P[j + 1] \) in \( P_Y \) in line 12 (this shrinks the range \([y_1^r, y_2^r] \)). Note that, once we know that we can descend from \( y_r \) by \( P[j + 1] \) (because \( r + \ell_r \geq j + 1 \)), we can compute the child node on the Patricia tree without accessing the text, both for explicit and virtual nodes \( y_r \). Thus, by computing \( \ell_r \) once when the active position \( r \) is created, in time \( O(\ell_r) \), we save all the accesses to \( T \) that would have been needed to descend from virtual nodes \( y_r \in P_Y \); when \( y_r \) is not the root, its text position is not phrase-aligned, so we cannot access its first symbols in constant time using \( A \).

Line 13 updates the nodes \( x_r \) of the surviving active positions, because some ranges \([x_1^r, x_2^r] \times [y_1^r, y_2^r] \) could be empty after we reduce \([y_1^r, y_2^r] \). For every active position \( r \), as long as there are no points in \([x_1^r, x_2^r] \times [y_1^r, y_2^r] \), we move \( x_r \) to its parent in \( P_X \). This process eventually terminates because, when \( x_r \) is the root and \([x_1^r, x_2^r] \) is the whole range of columns, we know that there are points in the band \([y_1^r, y_2^r] \) because it corresponds to the node \( y_r \).

Lines 8, 14, and 17 recompute the value \( p = \min \{ r - |x_r| + 1, r \text{ is active} \} \). This is necessary to make \( i \) grow as needed so that \( P[i..j + 1] \) occurs in \( T \), then reestablishing the invariant that \( P[i..j + 1] \) is the last MEM of \( P[i..j + 1] \). If \( p = i \), then \( P[i..j + 1] \) occurs in \( T \) (as it has a primary occurrence in some \([x_1^r, x_2^r] \times [y_1^r, y_2^r] \)), so we can retain the current value of \( i \); line 18 collects some text position \( t \) to be reported in case \( (i, j + 1) \) turns out to be a MEM of the whole \( P \). If, on the other hand, \( p > i \), this means \( P[i..j + 1] \) does not occur in \( T \) and thus \( (i, j) \) was a MEM. Lines 20–21 then report MEM \( (i, j) \) with its text position \( t \) (collected in the previous cycle of \( j \)) and increase \( i \) to \( p \), since only \( P[p..j + 1] \) occurs in \( T \). This could make \( i \) exceed \( j + 1 \) when the window becomes empty; otherwise line 23 finally inserts \( j + 1 \) as an active position. Line 26 reports the final MEM when \( j \) reaches \( m \).
Algorithm 2 Finding the MEMs of \(P[1 \ldots m]\) in \(T\) using a grammar-based index.

```plaintext
1  \(i \leftarrow 1; R \leftarrow \emptyset;\)
2  for \(j \leftarrow 0, \ldots, m - 1\) do
3      if \(i = j + 1\) and the root of \(P_X\) has no child labeled \(P[j + 1]\) then \(i \leftarrow i + 1;\)
4      else
5          \(y_{j+1} \leftarrow\) root of \(P_Y;\)
6          \(v \leftarrow\) descend in \(P_Y\) as much as possible with \(P[j + 2 \ldots]; \ell_{j+1} \leftarrow \|v\|;\)
7          \(x_{j+1} \leftarrow\) descend in \(P_X\) as much as possible with \(P[i \ldots j + 1]\)rev; \(r_{\text{min}} \leftarrow j + 1;\)
8
9      end for \(r \in R\) do
10     if \(r + \ell_r = j\) then \(R \leftarrow R \setminus \{r\};\)
11      else
12          \(y_r \leftarrow\) child of \(y_r\) by \(P[j + 1];\)
13          while \(r \neq \emptyset\) and \([x_r^1, x_r^2] \times [y_r^1, y_r^2]\) is empty do \(x_r \leftarrow\) parent of \(x_r;\)
14          if \(r - \|x_r\| < r_{\text{min}} - |x_{r_{\text{min}}}|\) then \(r_{\text{min}} \leftarrow r;\)
15      end
16  \(p \leftarrow r_{\text{min}} - |x_{r_{\text{min}}}| + 1;\)
17  if \(p = i\) then \(\ell \leftarrow\) text position of some point in \([x_r^1, x_r^2] \times [y_r^1, y_r^2]\); \(i \leftarrow p;\)
18  else
19      report \((i, j)\) with position \(T[t - j + i \ldots t];\)
20  end
21  if \(i \leq j + 1\) then \(R \leftarrow R \cup \{j + 1);\)
22 end
23 if \(i \leq m\) then report \((i, m)\) with position \(T[t - m + i \ldots t];\)
```

### 4.2 Analysis

For each value of \(j\), we spend \(O(1)\) time per active position. Since there are \(O(m)\) active positions at any time, this amounts to \(O(m^2)\) time.

The costs of lines 6, 7, and 13, are better charged to each active position \(r\), from its creation to its inactivation. When \(r\) is created, we spend \(O(m)\) time to compute \(\ell_r \leq m\) and \(x_r\) (since \(|x_r| \leq m\)). Later, we can decrease \(|x_r|\) several times, performing one range emptiness query in \([x_r^1, x_r^2] \times [y_r^1, y_r^2]\) per decrement of \(|x_r|\) (in fact we can go directly to the lowest physical ancestor of \(x_r\) rather than to its possibly virtual parent node, since otherwise the range \([x_r^1, x_r^2]\) will not change). Thus, we maintain overall \(O(m^2)\) emptiness queries, up to \(m\) per position \(r\) along its life. Maintaining the variables associated with active positions allows us amortizing these costs along the process.

Emptiness queries on \(G\) can be solved in \(O(\log g_{\epsilon})\) time and \(O(g_{\epsilon})\) space for any constant \(\epsilon > 0\) [9]; a recent construction takes \(O(g^{1/2})\) time [2]. The same complexity holds for returning one point in nonempty ranges. The \(O(m^2)\) cost charged to positions \(r\) is then multiplied by this factor. The rest of the construction time is inherited from the CFG-based index [12]; extending it to RLCFGs does not increase it.
Theorem 10. Let \( g_{t1} \) be the size of a RLCFG generating only \( T[1 \ldots n] \). Then, for any constant \( \epsilon > 0 \), we can build in \( O(g_{t1} \log^2 n) \) time a data structure of size \( O(g_{t1}) \) that finds the MEMs of any given pattern \( P[1 \ldots m] \) in time \( O(m^2 \log^2 g_{t1}) \subseteq O(m^2 (\log \delta + \log \log n)) \), with an occurrence of each. The query process uses \( O(m) \) additional space.

As mentioned, any CFG can also be used in the theorem. By using an emptiness structure of size \( O(g_{t1} \log \log g_{t1}) \) [9], we find the MEMs in time \( O(m^2 \log \log g_{t1}) \).

5 Indexing Locally Consistent Grammars

Before entering into the details of our more sophisticated solution, we must introduce some new concepts. A locally consistent grammar is a kind of RLCFG that guarantees that equal substrings of \( T \) are covered by similar subtrees of the parse tree, differing in \( O(1) \) nonterminals at each level of both subtrees. This has been used to produce grammar-based indices that find all the primary occurrences with only a logarithmic number of cuts in \( P \), thereby obtaining exact pattern searches in time that grows only linearly with \( m \) [11, 25, 24]. In this paper we make use of the latest result [24]. We present a lighter informal description; see the original paper for full details.

5.1 The Grammar

We first define the grammar [24, Sec. 3], which is produced level by level, for \( O(\log n) \) levels. Let \( S_k \) be the sequence of terminals and nonterminals forming level \( k \) of the grammar. Let \( \ell_k = (4/3)^{[k/2]} - 1 \), and let \( \mathcal{A}_k \) be the set of symbols \( A \) such that \( \exp(A) \leq \ell_k \). Those are the symbols that can be grouped to form new nonterminals in level \( k \).

Our string at level 0 is \( S_0 = T \). To form the string \( S_1 \), we detect the maximal runs of (at least 2) equal consecutive symbols in \( S_0 \) that are in \( \mathcal{A}_1 = \Sigma (\Sigma \) is the alphabet of \( T \) and also the set of terminals of the RLCFG). For each such run, say of \( t \) symbols \( a \in \mathcal{A}_1 \), we create the rule \( A \rightarrow a^t \) and replace the run by the nonterminal \( A \). The resulting sequence after all the runs have been replaced is \( S_1 = \text{rle}_{\mathcal{A}_1}(S_0) \), which contains terminals and nonterminals.

To form level 2, we define a function \( \pi_2 \) that reorders at random the distinct symbols of \( S_1 \), and use it to define blocks in \( S_1 \). Each position \( 0 < i < |S_1| \) such that

\[
\pi_2(S_1[i - 1]) > \pi_2(S_1[i]) < \pi_2(S_1[i + 1])
\]

is the end of a block. We also set ends of blocks at \( |S_1| \) and before and after every symbol not in \( \mathcal{A}_2 \) (which is still \( \Sigma \) per the formula of \( \ell_k \), so the runs introduced in \( S_1 \) cannot yet be grouped). For each distinct resulting block \( S_1[i \ldots j] \) we create a new rule \( A \rightarrow S_1[i \ldots j] \) and replace every occurrence of the same block in \( S_1 \) by \( A \). The resulting string is called \( S_2 = \text{bc}_{\pi_2, \mathcal{A}_2}(S_1) \). The process continues in the same way for odd and even levels:

\[
\begin{align*}
S_k &= \text{rle}_{\mathcal{A}_k}(S_{k-1}) \quad \text{if } k \text{ is odd}, \\
S_k &= \text{bc}_{\pi_k, \mathcal{A}_k}(S_{k-1}) \quad \text{if } k \text{ is even},
\end{align*}
\]

until we reach \( |S_k| = 1 \) for some \( k = O(\log n) \). The algorithm is Las Vegas type, trying out functions \( \pi_k \) to obtain some desired grammar size, but otherwise any functions \( \pi_k \) yield a correct index. They [24] prove that, in \( O(n) \) expected time, a RLCFG of size \( O(\delta \log \frac{n}{\ell}) \) is obtained, where \( \delta \) is a lower bound measure based on the substring complexity of \( T \) [11]: let \( T_\ell \) be the number of distinct length-\( \ell \) substrings in \( T \), then \( \delta = \max \{ T_\ell / \ell, \ell > 0 \} \). Interestingly,
for every $n$ and $\delta$, there exists a string family that requires $\Omega(\delta \log \frac{n}{\delta})$ space (i.e., $\log(n)$-bit words) to be represented [25]; therefore using space $O(\delta \log \frac{n}{\delta})$ for a grammar (and for an index) is asymptotically optimal for any specific $n$ and $\delta$.

A key property of this grammar is local consistency. Let $B_k$ be the set of all the ends of level-$k$ blocks:

$$B_k = \{\exp(S_k[i..j]), \ 1 \leq j \leq |S_k|\},$$

where we are extending $\exp(\cdot)$ homomorphically to strings. The cuts of level $k$ that fall inside the substring at $T[i..j]$ have the following positions inside $T[i..j]$:

$$B_k(i,j) = \{p - i + 1, \ p \in B_k \cap [i..j-1]\}.$$

Local consistency makes the sets $B_k(i,j)$ and $B_k(i',j')$ similar if $T[i..j] = T[i'..j']$, except at the extremes. Concretely, let $\alpha_k = [8\ell_k]$, then $B_k(i + 2\alpha_k, j - \alpha_k) = B_k(i' + 2\alpha_k, j' - \alpha_k)$.

An additional property of the resulting grammar is that it is locally balanced: the subtree of the parse tree rooted at nonterminal $A$ is of height $O(\log |\exp(A)|)$. This is a consequence of the fact that in $S_k$ there are fewer than $1 + 4(j - i + 1)/\ell_{k+1}$ blocks ending inside $T[i..j]$, and the height of $A$ is never more than the level $k$ of the string $S_k$ where it was created.

### 5.2 Pattern Searching

Let us now define which cuts of $P$ we need to try out in order to capture all the primary occurrences with this grammar [24, Sec. 4]. Since ends of blocks in $B_k(i,j)$ correspond to the phrase endings where a primary occurrence $T[i..j] = P$ can be cut, our set of cutting positions must suffice to capture those possible block endings for all $k$ and for every possible $T[i..j]$ that matches $P$. We define

$$M_k(i,j) = B_k(i,j) \setminus [2\alpha_{k+1} + 1..j - i - \alpha_{k+1}]$$

$$\cup \{\min(B_k(i,j) \cap [2\alpha_{k+1} + 1..j - i - \alpha_{k+1}])\},$$

that is, all the cutting points in the extremes, where the different occurrences of $T[i..j]$ may differ, and just the first one in the part that is guaranteed to be equal. Over all the levels,

$$M(i,j) = \bigcup_{k \geq 0} M_k(i,j).$$

The key point [24] is that $M(i,j)$ depends only on the content of $T[i..j]$ (not on its position in $T$), so we can define $M(P) = M(i,j)$ if $P = T[i..j]$, and this is the same set for every possible occurrence of $P$ in $T$. Further, $|M(P)| = O(\log m)$. In operational terms, this means that, at query time, we parse $P$ in $O(m)$ time using the same rules we defined for $T$, producing a parse tree of height $O(\log m)$ and finding the $O(\log m)$ cutting points $M(P)$.

### 6 A Faster Solution using Locally Consistent Grammars

The idea to use the index of the preceding section is to exploit the fact that $O(\log(j-i+1))$ cutting points suffice to find all the primary occurrences of any window $P[i..j]$. We will then maintain the parse tree of $P[i..j]$, and the set $M(P[i..j])$, as we slide it through $P$, and use it to maintain the number $|R|$ of active positions within $O(\log m)$. We also need more sophisticated mechanisms to avoid the quadratic costs in lines 6, 7, and 13 of Algorithm 2.
6.1 Parsing the Pattern

In this section we show how we maintain the parse tree of $P[i...j]$, or more precisely, the corresponding strings $S_0, S_1, \ldots$, as well as $M(P[i...j])$ and $R$, as we slide $P[i...j]$ along $P$. Recall that the height of the parse tree of $P$ is $O(\log m)$ in our locally-balanced grammar.

Maintaining the parse

Assume the parse tree is built for $P[i...j]$ and now we have to increment $j$. At level $k = 0$, we simply extend $S_0$ by the symbol $e_0 = P[j + 1]$. This propagates upwards as follows, where $l_k$ is the last symbol of $S_k$ and $e_{k-1}$ has just been added at the end of $S_{k-1}$.

1. If $k$ is odd (a run-formation level), $l_k = l_{k-1}$ or $l_k \rightarrow l_{k-1}^t$, $|\text{exp}(e_{k-1})| \leq \ell_k$, and $l_{k-1} = e_{k-1}$, we find or create a rule $e_k \rightarrow l_{k-1}^{t+1}$ ($t = 1$ if $l_k = l_{k-1}$) and replace $l_k$ by $e_k$.

2. If $k$ is even (a block-formation level), $l_k = l_{k-1}$ or $l_k \rightarrow \beta l_{k-1}$, $|\text{exp}(e_{k-1})| \leq \ell_k$, and $\pi_k(l_{k-1}) > \pi_k(e_{k-1})$, we find or create a rule $e_k \rightarrow \beta l_{k-1} e_{k-1}$ ($\beta = \varepsilon$ if $l_k = l_{k-1}$) and replace $l_k$ by $e_k$.

3. In any other case, we just append $e_k = e_{k-1}$ at the end of $S_k$.

We see that insertions at the end of $S_{k-1}$ propagate as new insertions or replacements at the end of $S_k$. We process those replacements as the deletion of $l_k$ followed by the insertion of $e_k$ at the end. The following are the rules to propagate to $S_k$ the deletion of $l_k$.

1. If $l_k = l_{k-1}$, we delete $l_k$.

2. If $l_k \rightarrow l_{k-1}^t$ is a run, we find or create the rule $e_k \rightarrow l_{k-1}^{t-1}$ (just $e_k = l_{k-1}$ if $t - 1 = 1$) and replace $l_k$ by $e_k$.

3. If $l_k \rightarrow \beta l_{k-1}$ is a block, we find or create the rule $e_k \rightarrow \beta$ (just $e_k = \beta$ if $|\beta| = 1$) and replace $l_k$ by $e_k$.

Each of those updates can be carried out in constant time by just maintaining linked lists with the sequence of symbols in each string $S_k$, perfect hash tables with the existing right-hand sides of the run-formation rules $l_k \rightarrow l_{k-1}^t$, and tries with the right-hand sides of the block-formation rules $l_k \rightarrow \beta l_{k-1}$. In particular, each block-formation nonterminal $l_k$ points to the node in the trie that represents its string $\beta l_{k-1}$ and the trie node representing $\beta l_{k-1}$ stores $l_k$. With parent pointers in the trie, we have constant-time access to the node of $\beta$ from the node of $\beta l_{k-1}$, and with child pointers we move from $\beta l_{k-1}$ to $\beta l_{k-1} e_{k-1}$. The children of trie nodes are stored in perfect hash tables to enable downward traversals in constant time. All this can be precomputed in expected time linear in the grammar size.

The case when $i$ increases is symmetric. We start by deleting the first symbol of $S_0$, and propagate the update upwards acting on the first symbols $f_k$ at each level $k$. To handle those operations we need the lists for $S_k$ be doubly-linked, and also to store tries for all the right-hand sides read as $f_k \rightarrow f_{k-1}^\prime$.

The number of updates are actually bounded to $O(1)$ updates per level, and thus to $O(\log m)$ per increase of $j$ or of $i$. Consider the string $P[i...j].S$, where $S$ is a special symbol for which we assume $\pi_k(S) = +\infty$ for all $k$. The parse tree of $P[i...j].S$ is then identical to that of $P[i...j]$, just adding $S$ at the end of every $S_k$. The strings $S_k$ formed for $P[i...j]$ followed by $S$ are the same formed for $P[i...j]$ followed by $P[j + 1]$, except for the first $2\alpha_k$ and the last $\alpha_k$ positions of $P[i...j]$ [24, Lem. 3.7]. Therefore, the addition of $P[j + 1]$ can only alter the last $1 + \alpha_k$ positions of $P[i...j + 1]$ in each $S_k$. Analogously, removing $P[i]$ can only alter the first $2\alpha_k - 1$ positions of $P[i + 1...j]$ in its strings $S_k$. On the other hand, those $O(\alpha_k)$ positions correspond to only $O(1)$ symbols in $S_k$ [24, Lem. 3.8]. The total amount of work is proportional to the number of updated symbols if we perform them levelwise, on $S_0$, then on $S_1$, and so on. All the changes then add up to $O(m \log m)$ along the processing of $P$. 

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24:10 Computing MEMs on Repetitive Text Collections
Dealing with unknown symbols

We analyzed the parsing process as if we would always find a known nonterminal for the
right-hand sides we modify, but it could be that we have to create new nonterminals that
were never formed during the parsing of $T$.

To handle those cases, we create fresh nonterminals and continue the process normally,
removing them when they are no longer needed. An easy way to handle this would be to
make the hash tables and tries of right-hand sides dynamic, so we can add and remove
elements in the tables and nodes; we will soon sharpen this solution.

We must assign values $\pi_k(e_k)$ to the fresh symbols $e_k$ we create in $S_k$. We can assign
arbitrary values (different from the other $\pi_k$ values) without affecting correctness: the index
works correctly for arbitrary functions $\pi_k$, as explained. No matter how we choose the values
$\pi_k(e_k)$, we can add $O(m \log m)$ fresh symbols along the whole process, but we can do better.

New symbols $e_k$ may appear in the parsing of $P[i..j]$ even if $P[i..j]$ appears in $T$,
because the parsing of $P[i..j]$ can be different from that of its occurrences in $T$. However,
this can happen only in the first $2\alpha_k$ and the last $\alpha_k$ positions, in $S_k$. Once the end of $e_k$
falls before position $j - \alpha_k$, and it is after the position $2\alpha_k$, then $e_k$ should have appeared in
the parsing of every occurrence of $P[i..j]$ in $T$ [24, Lem 3.7]. Therefore, when we completely
incorporate $P[j+1]$ and as a result the end of a fresh symbol $e_k$ of $S_k$ falls behind position
$j+1-\alpha_k$ of $P[i..j+1]$, we know $P[i..j+1]$ cannot appear in $T$ until the position falls
behind $i+2\alpha_k$. At this point, then, we can suspend the search (very much as Algorithm 2
does in line 3) and increase $i$ until $e_k$ ends within the leftmost $2\alpha_k$ symbols of $P[i..j+1]$.

This has as a consequence that we can have only $O(1)$ fresh symbols per level, just as
$M_{i,j}(P)$, and $O(\log m)$ in total. Instead of making the tries and hash tables dynamic, we
can have one extra atomic heap per hash table (the one for the run-length symbols and the
one in each trie node) where we can insert/delete the necessary fresh symbols, and they will
be processed in constant time. We then retain the $O(m \log m)$ total processing cost.

Maintaining $M(P)$ and $R$

After we finish updating the parse tree, we collect the first $2\alpha_{k+1}$ positions, the position of the
following end of block, and the last $\alpha_{k+1}$ positions, in each list $S_k$, to form the sets $M_k(P)$.
Those are then merged into $M(P)$ and sorted by increasing value. Since $|M(P)|=O(\log m)$,
and its values are integers in $[1..m]$, $M(P)$ can be sorted in $O(\log m)$ time with atomic
heaps. We then traverse $M(P)$ and the current set $R$ in synchronization, so as to (1) remove
the positions of $R$ that are not anymore in $M(P)$, and (2) insert in $R$ the positions that are
now in $M(P)$, as long as the position had not been in $R$ before and had been inactivated
(this is easily marked in an array of size $m$). At the end of this process, it always holds that
$R \subseteq M(P[i..j])$, and thus $|R|=O(\log m)$. Due to the parsing, an active position $r$ may
enter and leave $R$ several times along the process, but this time that will not be an issue.

Overall, we maintain the parsing, $M(P[i..j])$, and $R$ in time $O(m \log m)$. Since all the
lines in Algorithm 2 other than 6, 7, and 13, take $O(1)$ time per element in $R$, the total time
spent in those lines adds up to $O(m \log m)$.

6.2 Patricia Tree Searches

Lines 6 and 7 of Algorithm 2 perform $\Theta(m)$-time searches in $P_X$ and $P_Y$. Since each of the
$m$ positions becomes active when $j+1$ reaches it, this amortizes to no less than $\Theta(m^2)$,
which is now too high for us. We then resort to a different technique.
Instead of computing $\ell_{j+1}$ and $x_{j+1}$ inside the main cycle, we will compute them all beforehand, in batch form. We make use of the following result, which was key to obtain subquadratic times in grammar-based indexing.

**Lemma 11** ([11, Lem 6.5]). Let $S$ be a set of strings and assume we have a data structure supporting extraction of any length-$t$ prefix of strings in $S$ in time $f_h(\ell)$ and computation of a given Karp–Rabin signature $\kappa$ of any length-$\ell$ prefix of strings in $S$ in time $f_h(\ell)$. We can then build a data structure of $O(|S|)$ words such that, later, given a pattern $P[1..m]$ and $\tau$ suffixes $Q_1, \ldots, Q_\tau$ of $P$, we find the ranges of strings in (the lexicographically-sorted) set $S$ prefixed by each $Q_i$, in $O(m + \tau(f_h(m) + \log m) + f_L(m))$ total time.

When $S = \mathcal{X}$ or $S = \mathcal{Y}$, our access data structure $A$ provides the required prefix/suffix extraction in time $f_h(\ell) = O(\ell)$. As for Karp-Rabin signatures [20], a result of independent interest is that we can obtain $f_h(\ell) = O(\log \ell)$ time on our grammar, as proved next. We consider the more complicated case of $Y \in \mathcal{Y}$; the case of $X \in \mathcal{X}$ is analogous. Recall we can compute in $O(1)$ time one of $\kappa(S \cdot S')$, $\kappa(S)$, and $\kappa(S')$, from the other two [11, Sec. A.3].

**Lemma 12.** The Karp-Rabin signature $\kappa(Y[1..\ell])$ of any $Y \in \mathcal{Y}$ can be computed in time $O(\log^2 \ell)$ with our grammar.

**Proof.** We build on the same structure $A$ used for extraction from the root of $P$. The strings in $\mathcal{Y}$ are concatenations $Y = \exp(B_s_1) \cdots \exp(B_i)$ of siblings in rules $A \rightarrow B_1 \cdots B_i$ in the grammar tree. The node $v \in P_\mathcal{Y}$ of $Y$ stores $\langle v \rangle = B_s$. Let us first assume that $|\exp(B_s)| \geq \ell$, so the signature can be computed on $\exp(B_s)[..\ell]$.

Structure $A$ is a set of tries on the grammar symbols. The terminals $\Sigma$ form the trie roots. If $A \rightarrow B_1 \cdots B_i$, then $B_1$ is the parent of $A$. If $A \rightarrow B'$, then $B$ is the parent of $A$. Any ancestor $C$ of $B_s$ in the tries is a node that descends from $B_s$ by the leftmost path in the parse tree. The structure $A$ can jump from $B_s$ to any such $C$ in constant time in the tries. Our grammar is locally balanced: there can be only one block ending inside $\exp(B_s[..\ell])$ at level $k = 1 + 2 \log_{4/3}(4\ell)$ [24, Lem. 3.8], and thus the lowest $C$ such that $|\exp(C)| \geq \ell$ has level at most $k + 1$; its height is $d \leq k + 1 = O(\log \ell)$. It can then be found in $O(\log \log \ell)$ time with exponential search on the ancestors of $B_s$. We then have that $\exp(B_s)[..\ell] = \exp(C)[..\ell]$ and can compute the signature on $C$ instead.

The basic algorithm to compute signatures takes time $O(\log^2 \ell)$ [11, Lem 6.7]. It moves from $C$ towards the leaf $L$ of the parse tree that corresponds to $\exp(C)[\ell]$. Let $C \rightarrow C_1 \cdots C_l$, then it stores every $w_i(C) = |\exp(C_1 \cdots C_l)|$ and every $\kappa_i(C) = \kappa(\exp(C_1 \cdots C_l))$. The algorithm finds, in $O(\log \ell)$ time, using exponential search, the $C_i$ that is in the path to $L$ (i.e., $w_{i-1} < \ell \leq w_i$), sets $\ell \leftarrow \ell - w_{i-1}$, collects $\kappa_{i-1}(C)$, and continues by $C_i$. It composes all these $\kappa$ values towards $L$ to obtain $\kappa(Y[1..\ell])$. In rules $C \rightarrow C^*_l$ it obtains $i$ in constant time but spends $O(\log i)$ time to compute $\kappa_{i-1}(C)$ from the stored $\kappa(\exp(C_l))$.

Instead, an $O(\log n)$ time algorithm [11, Thm. A.3] replaces the exponential searches by a more sophisticated scheme, whose cost is the telescoping sum $\sum_{h=1}^{l} \log(t_h/t_{h-1}) \leq \log t_p$, where $t_h$ is the number of children (counting $C \rightarrow C^*_l$ as having $t$ children) of the ancestor at distance $h$ of leaf $L$. In their case, they start from the root, which could have $t_p = n$, but if we start it from a node $C$, its time is $\log t_p \leq \log |\exp(C)|$. Another component of the cost is the number of times one leaves from heavy paths; this is again $O(\log n)$ in general but just $O(d) = O(\log \ell)$ if we start from the position of $C$ in its heavy path.

It could be, however, that $|\exp(C)|$ is as long as $n$. Because it was formed in $S_k$, however, the children $C_i$ of $C$ belonged to $A_k$ (only those nonterminals are allowed to form rules in $S_k$), and thus by definition $|\exp(C_i)| \leq \ell_k$ and $\log |\exp(C_i)| = O(k) = O(\log \ell)$. We can then
find $i$ and compute $\kappa_{i-1}(C)$ in time $O(\log i) \subseteq O(\log \ell)$ with the basic method \cite[Lem 6.7]{G. Navarro 24:13} and then continue from $C_i$, where the more sophisticated technique \cite[Thm. A.3]{G. Navarro 24:13} completes the computation in another $O(\log |exp(C_i)|) \subseteq O(\log \ell)$ time.

In case $|exp(B_s)| < \ell$, we find the first $s < i \leq t$ such that $w_i(A) \geq \ell$, and compute instead the signature of $exp(B_s)\ldots \ell - w_{i-1}(A)]$, to then compose it with the stored values $\kappa_{s-1}(A)$ and $\kappa_{i-1}(A)$ to obtain the final signature $\kappa(Y[\ldots \ell]) = \kappa(exp(A)[w_{s-1}(A) + 1 \ldots w_{s-1}(A) + \ell])$.

### Batched searches

The $m$ searches for all the values $\ell_r$, $1 \leq r \leq m$, correspond to searching $P_Y$ for every suffix $P[r+1\ldots]$. Note that Lemma 11 does not yield the node $v$ of line 6, but rather its corresponding range $[v^1, v^2]$. By performing a lowest common ancestor (LCA) query on $P_Y$ from the $v^1$th and $v^2$th leaves, we obtain $v = lca(v^1, v^2)$ (identifying leaves with their ranks). The answer is indeed $v$ if $|v| = m - r$; if $m - r < |v|$ the answer is the virtual node of string length $m - r$ on the edge of $P_Y$ that leads to $v$. Linear-space LCA data structures that are built in linear time and answer $lca$ in $O(1)$ time are well known \cite{G. Navarro 24:13}.

The problem is that Lemma 11 works only if $P[r+1\ldots]$ actually prefixes some string in $Y$. Otherwise, unlike classical trie searching, it does not even yield the maximum prefix of $P[r+1\ldots]$ that prefixes some string in $Y$. We will resort to, essentially, binary searching for those longest prefixes using Lemma 11 as an internal tool.

Assume $m$ is a power of 2 for simplicity; the general case is easily deduced. We define sets $Q_{a,b}$ of positions, containing those values $r$ such that $P[r+1\ldots a]$ is known to be a prefix in $Y$ and $P[r+1\ldots a+b+1]$ is known not to be a prefix in $Y$ (the first condition is assumed to hold if $r+1 > a$). We start with the set $Q_{1,m} = \{1, \ldots, m\}$. To process a set $Q_{a,b}$, we search for all the $\tau = |Q_{a,b}|$ suffixes $\{P[r+1\ldots c] \mid r \in Q_{a,b}\}$ of $P[\ldots c]$ using Lemma 11, with $c = (a + b + 1)/2$. The values $r$ where $P[r+1\ldots c]$ is found are moved to $Q_{c,b}$, and the others to $Q_{a,c-1}$ (if $r + 1 > c$, then $P[r+1\ldots c] = \varepsilon$, so we can directly move $r$ to $Q_{c,b}$ without searching for it). We will associate the node $v_{r,c} \in P_Y$ to those values $r$ for which $P[r+1\ldots c]$ is found in $Y$; those not found retain their previous node $v_{r,*}$ (in the beginning all such nodes are $v_{r,r}$ and equal to the root of $P_Y$).

Note that the values $b - a + 1$ halve as the elements in $Q_{a,b}$ are separated into two sets. Any value $r$ is then moved $O(\log m)$ times until it ends up in a set of the form $Q_{a,c}$; at this point we know that the longest suffix of $P[r+1\ldots]$ that is also a prefix in $Y$ is $P[r+1\ldots c]$, and also know its node $v_{r,c}$.

The cost of using Lemma 11 has two parts. The cost $f_a(m) + \log m = O(\log m)$ can be charged to each of the $\tau$ suffixes sought, and there is an additional global cost of $m + f_a(m) = O(m)$. Since every suffix $P[r+1\ldots]$ participates $O(\log m)$ times in the lemma, the first cost adds up to $O(m \log^2 m)$ over all the $m$ positions $r$. The second part is potentially very large, however: the suffixes in $Q_{a,b}$ may start well ahead of $a$, thus the pattern is $P[\ldots c]$, not $P[a\ldots c]$; a simple application of the lemma would lead to a quadratic cost again.

### Smarter substring extractions

To reduce this time, we consider where the $O(m)$ cost in Lemma 11 comes from. A first part refers to the time needed to compute the Karp-Rabin signatures for all the suffixes in $Q_{a,b}$. This cost is easily maintained within $O(m)$ overall because we can compute the signatures $\kappa(P[r+1\ldots])$, for all $1 \leq r \leq m$, in a single pass over $P$, and then any $\kappa(P[r+1\ldots j])$ is obtained in constant time from $\kappa(P[r+1\ldots])$ and $\kappa(P[j+1\ldots])$. 


The second part of the $O(m)$ cost corresponds to the time $f_e(m)$ to verify the longest suffix among those that passed some previous filters; the rest of the verification is built on that extracted suffix. Let $P[r+1..c]$ be the longest candidate suffix. If $r+1 > a$, we extract the actual suffix $P'[..c-r]$ regularly in time $f_e(c-r) = O(b-a)$ with $A$, because $P'$ starts at the root of $P_Y$ and thus it belongs to $\mathcal{Y}$.

Otherwise, $r+1 \leq a$ and thus $P[r+1..a]$ had been successfully matched before and we have its node $v_{r,a} \in P_Y$. As mentioned, the process of Lemma 11 performs several checks before performing the final extraction of the longest suffix surviving the checks. We will add a new check to those, which can only speed up the process: the candidate node $v$ for $P[r+1..c]$ must now descend from $v_{r,a}$ in order to be further considered. The descendant check is performed in constant time by comparing the leaf range $[v^1,v^2]$ of $v$ with that of $v_{r,a}$. If $v$ passes the test, we know that $P'[..c-r]$ does start with $P[r+1..a]$, and then only need to extract $P'[a-r+1..c-r]$ from the text, which is of length $O(b-a)$.

This time, however, the string to extract does not start at the root of $P_Y$, and thus it requires a random access to $T$. Recall, as in Lemma 12, that the strings in $\mathcal{Y}$ are concatenations $Y = \text{exp}(B_1) \cdots \text{exp}(B_t)$ of consecutive siblings in rules $A \rightarrow B_1 \cdots B_t$ in the grammar tree (if $A \rightarrow B'$, then the node stores $(v) = B^{(t-1)}$ and we have $B_1 = B$). Let us first assume that $|\text{exp}(B_k)| \geq c$, so the substring to extract is within $\text{exp}(B_k)[..c]$. We use again the structure $A$, now to extract the string in time $O(b-a + \log c)$.

Once again, we can search in $O(\log \log c)$ time for the lowest descendant $C$ of $B_k$ such that $|\text{exp}(C)| \geq c$; its height is $d = O(\log c)$ because the grammar is locally balanced. Since $\text{exp}(B_k)[..c] = \text{exp}(C)[..c]$, we descend from $C$ to the leaf $L$ in the parse tree representing $\text{exp}(C)[a-r+1]$. Using the same techniques as in Lemma 12, the time is $O(d) = O(\log c)$.

From $L$, $\text{exp}(C)[a-r+1..c-r] = P'[a-r+1..c-r]$ is extracted in time $O(c-a) = O(b-a)$.

In case $|\text{exp}(B_k)| < c$, the node $C$ is not a descendant of $B_k$ but we use $C = A$ instead. Given the limitation on $|\text{exp}(B_k)|$, the height of $B_k$ is $O(\log c)$, and so is the height of $A$.

The $O(\log c) = O(\log m)$ cost can be charged to the suffix sought, which adds up to $O(m \log^2 m)$ over the $O(\log m)$ times each suffix may use the lemma. The $O(b-a)$ terms add up to $O(m)$ per level of sets $Q_{a,b}$ (a level corresponding to a difference $b-a+1$). Since all the ranges $(a,b)$ of a level are disjoint (level $\ell$ partitions $(1,m)$ into $2^\ell$ ranges of size $m/2^\ell$), the $b-a$ values add up to $O(m)$ per level. Since there are $O(\log m)$ levels, that part of the cost adds up to $O(m \log m)$.

We similarly compute the nodes $x_r$ for every $P[..r]^{rev}$ on $P_Y$. While in line 7 of Algorithm 2 we search only for $P[i..r]^{rev}$ because we are not interested in positions before $i$, this time we precompute all the values in advance for the smallest $i = 1$. Later, when $i$ increases, we will move to the required ancestors of $x_r$ in line 13.

Overall, the Patricia trie searches execute lines 6 and 7 in $O(m \log^2 m)$ total time.

### 6.3 Emptiness Queries

In line 13, we perform one range emptiness query every time we decrement $|x_r|$ for some $r$; this amounts to $O(m^2)$ emptiness queries, which we cannot afford now. We will instead use a faster method based on orthogonal range successor queries: given a range $[x^1_r, x^2_r] \times [y^1_r, y^2_r]$,
we can find the largest value \(x_\leq \leq x_1\) such that \([x_\leq, x_1]\times[y_1, y^2]\) contains a point, and the smallest value \(x_\geq \geq x_2\) such that \([x_1, x_\geq]\times[y_1, y^2]\) contains a point. Those queries can run in \(O(\log^\epsilon g)\) time on a grid with \(g\) points, using an \(O(g)\)-space data structure, for any constant \(\epsilon > 0\) defined at construction [35]; construction time can be made \(O(g^{\sqrt{\log g}})\) [2].

The lowest ancestor \(x\) of \(x_i\) containing some point in \([x_1, x^2]\times[y_1, y^2]\) must then hold \(x_1 \leq x_i \leq x_2\). In the first case, it is \(v_1 = \text{lca}(x_\leq, x_2)\); in the second, it is \(v_2 = \text{lca}(x_1, x_\geq)\). Both \(v_1\) and \(v_2\) are ancestors of \(x\), and thus of each other. The correct node \(x\) is then the lowest of \(v_1\) and \(v_2\), which is known from the leaf ranges stored at the nodes.

With this query, line 17 of Algorithm 2 does not cycle; it just performs one iteration step. It can then be counted as one of the \(O(|R|)\) operations performed in each cycle \(j\). Since there are \(O(m \log m)\) such operations, this one adds \(O(m \log m \log^\epsilon g)\) to the total cost.

### 6.4 The Final Result

Our time complexities then add up to \(O(m \log m (\log m + \log^\epsilon g))\) for a grammar of size \(g\). Since in our case \(g = O(\delta \log \frac{n}{\delta})\), we can write the time as \(O(m \log m (\log m + \log^\epsilon \delta + \log \log n))\). The construction time of all the data structures we use is dominated by the \(O(n \log n)\) expected time needed to build the Karp-Rabin hashes of Lemma 11 [11, Sec. 6.6] (the grammar is built in \(O(n)\) expected time, see [24, Cor. 3.15]).

**Theorem 13.** Let \(T[1..n]\) have substring complexity \(\delta\). Then, for any constant \(\epsilon > 0\), we can build in \(O(n \log n)\) expected time a data structure of size \(O(\delta \log \frac{n}{\delta})\) that finds the MEMs of any given pattern \(P[1..m]\) in time \(O(m \log m (\log m + \log^\epsilon \delta + \log \log n))\) \(\subseteq O(m \log m (\log m + \log^\epsilon n))\), with an occurrence of each. The query process uses \(O(m)\) additional space.

We have assumed \(m \leq n\), but it could be the other way in some applications. Since in this case no substring longer than \(n\) will be matched inside \(P\), we can run \(O(m/n)\) iterations finding the MEMs of \(P[1..2n], P[n..3n], P[2n..4n]\), and so on, avoiding repeated MEMs in the output. The total cost would then be \(O(m \log^2 n)\) and the query space would be \(O(n)\).

### 7 Conclusions

We have obtained improved results, including the first subquadratic algorithm, to find MEMs on parsing-based indices, which are the most promising in terms of space for highly repetitive text collections. While suffix-based indices can preprocess \(T[1..n]\) to find the MEMs of \(P[1..m]\) in \(T\) in time \(O(m \log \log n)\), their space is \(\Omega(r)\), where \(r\) (the number of runs in the BWT of \(T\)) is not such a strong measure of repetitiveness [34]. Our first result is a data structure of size \(O(g_{rl})\), where \(g_{rl}\) is the size of the smallest RLCFG that generates \(T\). This is currently the best possible space for any structure able to access \(T\) with relevant time guarantees [34]. Our structure finds the MEMs in \(O(m^2 \log^\epsilon n)\) time for any constant \(\epsilon > 0\). This is very similar to the time of previous work [18], which could also run in \(O(g_{rl})\) space. Within \(O(\delta \log \frac{n}{\delta})\) space, we obtain the first subquadratic time, \(O(m \log m (\log m + \log^\epsilon n))\), on a particular RLCFG that has local consistency properties. This space is optimal for every \(n\) and \(\delta\), though \(g_{rl}\) is always \(O(\delta \log \frac{n}{\delta})\) and can be \(o(\delta \log \frac{n}{\delta})\) in some text families [25].

A challenge for future work is to extend our results to finding \(k\)-MEMs, which are the maximal substrings of \(P\) that appear at least \(k\) times in \(T\), for \(k\) given at query time. The basic Algorithm 1 is easily modified to find the \(k\)-MEMs in \(O(m)\) time, but running in compressed space is more costly. In the extended version we will show how find \(k\)-MEMs, changing the \(\log^\epsilon n\) terms to \(\log^{2+\epsilon} n\), by building on grammar-based indices that can count
the number of occurrences associated with a set of primary occurrences \[11\]. The space also increases: the \(O(g_{\text{rl}})\) space becomes \(O(g)\) \((g \geq g_{\text{rl}}\) being size of the smallest CFG) and the \(O(\delta \log \frac{n}{\delta})\) space becomes \(O(\gamma \log \frac{n}{\gamma})\) \((\gamma \geq \delta\) being the size of a string attractor of \(T\) \[22\]).

Our techniques are presented on a particular locally consistent grammar \[24\] that yields the best complexities, but they would work on others too. We plan to implement them on practical constructions of CFGs \[12\] built with RePair \[27\] or of locally consistent grammars based on induced suffix sorting \[14, 36\]. Further, even without having theoretical guarantees, the algorithm for arbitrary RLCFGs will probably be competitive if implemented on Lempel-Ziv based indices \[26, 16\], which are considerably smaller than those based on grammars.

References


Abstract

In order to use them for compression, we extend L-systems (without ε-rules) with two parameters \(d\) and \(n\), and also a coding \(\tau\), which determines unambiguously a string \(w = \tau(\varphi^d(s))[1:n]\), where \(\varphi\) is the morphism of the system, and \(s\) is its axiom. The length of the shortest description of an L-system generating \(w\) is known as \(\ell\), and it is arguably a relevant measure of repetitiveness that builds on the self-similarities that arise in the sequence.

In this paper, we deepen the study of the measure \(\ell\) and its relation with a better-established measure called \(\delta\), which builds on substring complexity. Our results show that \(\ell\) and \(\delta\) are largely orthogonal, in the sense that one can be much larger than the other, depending on the case. This suggests that both mechanisms capture different kinds of regularities related to repetitiveness.

We then show that the recently introduced NU-systems, which combine the capabilities of L-systems with bidirectional macro schemes, can be asymptotically strictly smaller than both mechanisms for the same fixed string family, which makes the size \(\nu\) of the smallest NU-system the unique smallest reachable repetitiveness measure to date. We conclude that in order to achieve better compression, we should combine morphism substitution with copy-paste mechanisms.

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1 Introduction

In areas like Bioinformatics, it is often necessary to handle big collections of highly repetitive data. For example, two human genomes share 99.9% of their content [23]. In another scenario, for sequencing a genome, one extracts so-called reads (short substrings) from it, with a “coverage” of up to 100X, which means that each position appears on average in 100 reads.¹ There is a need in science and industry to maintain those huge string collections in compressed form. Traditional compressors based exclusively on Shannon’s entropy are not good for handling repetitive data, as they only exploit symbol frequencies when compressing. Finding good measures of repetitiveness and also compressors exploiting this repetitiveness has then become a relevant research problem.

¹ https://www.illumina.com/science/technology/next-generation-sequencing/plan-experiments/coverage.html
A strong theoretical measure of string repetitiveness introduced by Kociumaka et al. [12] is $\delta$, based on the substring complexity function. This measure has several nice properties: it is computable in linear time, monotone, resistant to string edits, insensitive to simple string transformations, and it lower-bounds almost every other theoretical or ad-hoc repetitiveness measure considered in the literature. Further, although $O(\delta)$ space is unreachable, there exist $O(\delta \log(n/\delta))$-space representations supporting efficient pattern matching queries [12, 11], and this space is tight: no $o(\delta \log(n/\delta))$-space representation can exist [12].

The idea that $\delta$ is a sound lower bound for repetitiveness is reinforced by the fact that it is always $O(b)$, where $b$ is the size of the smallest bidirectional macro scheme generating a string [26]. Those macro schemes arguably capture every possible way of exploiting copy-paste regularities in the sequences. Some very recent works [19], however, explore other sources of repetitiveness, in particular self-similarity, and are shown to break the lower bound of $\delta$.

The simplest of those schemes, which reuse the name L-system for simplicity [19], builds upon Lindenmayer systems [15, 16], in particular on the variant called CPD0L-systems. A CPD0L-system describes the language of the images, under a coding $\tau$, of the powers of a non-erasing morphism $\varphi$ starting from an string $s$ (called the axiom), that is, the set $\{\tau(\varphi^i(s)) \mid i \geq 0\}$. The L-systems extend CPD0L-systems with two parameters $d$ and $n$, so as to unambiguously determine the string $w = \tau(\varphi^d(s))[1 : n]$. The size of the shortest description of an L-system generating $w$ in this fashion is called $\ell$. Intuitively, $\ell$ works as a repetitiveness measure because any occurrence of the symbol $a$ at level $i$ expands to the same string at level $i + j$ for every $j$.

Since $\ell$ is a reachable measure of repetitiveness (because the L-system is a representation of $w$ of size $O(\ell)$), there are string families where $\delta = o(\ell)$. Intriguingly, it has been shown [19] that there are other string families where $\ell = o(\delta)$, so (1) both measures are not comparable and (2) the lower bound $\delta$ does not capture this kind of repetitiveness. On the other hand, it is shown that $\ell = O(g)$, where $g$ is the size of the smallest deterministic context-free grammar generating only $w$. This comparison is relevant because L-systems are similar to grammars, differing in that they have no terminal symbols, so their expansion must be explicitly stopped at level $d$ and then possibly converted to terminals with $\tau$.

Grammars provide an upper bound to repetitiveness that is associated with well-known compressors, so this upper bound makes $\ell$ a good measure of repetitiveness.

A more complex scheme that was also introduced [19] are NU-systems, which combine the power of L-systems with bidirectional macro schemes. The measure $\nu$, defined as the size of the smallest NU-system generating $w$, naturally lower bounds both $\ell$ and $b$. The authors could not, however, find string families where $\nu$ is asymptotically better than both $\ell$ and $b$, so it was unclear if NU-systems are actually better than just the union of both underlying schemes.

In this paper we deepen the study of the relations between these new intriguing measures and more established ones like $\delta$ and $g$. Our results are as follows:

1. We show that $\ell$ can be much smaller than $\delta$, by up to a $\sqrt{n}$ factor, improving a previous result [19] and refuting their conjecture that $\ell = \Omega(\delta/\log n)$.

2. On the other hand, we expose string families where $\ell$ is larger than the output of several repetitiveness-aware compressors like the size $g_{\text{rl}}$ of the smallest run-length context-free grammar, the size $z_e$ of the greedy LZ-End parse [13], and the number of runs $r$ in the Burrows-Wheeler Transform of the string [2]. We then conclude that $\ell$ is incomparable to almost all measures other than $g$, which suggests that the source of repetitiveness it captures is largely orthogonal to the typical cut-and-paste of macro schemes.
An alphabet $\Sigma$ is a finite set of symbols and is usually denoted by $\Sigma$. A (finite) string $w$ is a finite sequence $w[1]w[2] \ldots w[n]$ of symbols where $w[i] \in \Sigma$ for $i \in [1,n]$, and its length is denoted by $|w| = n$. The empty string, whose length is 0, is denoted by $\varepsilon$. The set of all possible finite strings over $\Sigma$ is denoted by $\Sigma^*$. If $x = x[1] \ldots x[n]$ and $y = y[1] \ldots y[m]$, the concatenation operation $x \cdot y$ (or just $xy$) yields the string $x[1] \ldots x[n]y[1] \ldots y[m]$. Let $w = xyz$. Then $y$ (resp. $x, z$) is a substring (resp. prefix, suffix) of $w$. It is proper if it is not equal to $w$, and non-trivial if it is distinct from $\varepsilon$ and $w$. The notation $w[i : j]$ stands for the substring $w[i] \ldots w[j]$ if $i \leq j$, and $\varepsilon$ otherwise. We also use the conventions $w[i : j] = w[1 : j]$ if $i < 1$, $w[i : j] = w[i : n]$ if $j > n$, and $w[i : j] = \varepsilon$ if $i > n$ or $j < 1$. Other convenient notations are $w[: i] = w[1 : i]$ and $w[: i] = w[i : |w|]$ for prefixes and suffixes, respectively.

A (right) infinite string $w$ (we use boldface to emphasize them) over an alphabet $\Sigma$ is a mapping from $\mathbb{Z}^+$ to $\Sigma$, and its length is called $\omega$, which is greater than $n$ for any $n \in \mathbb{Z}^+$. It is possible to define the concatenation $x \cdot y$ if $x$ is finite and $y$ infinite. The concepts of substring, prefix, and suffix carry over to infinite strings, with proper prefixes always being finite and suffixes always being infinite. The notations $w[i], w[i : j], w[: i]$ and $w[i :] = w[i]w[i+1] \ldots$ also carry over to infinite strings.

3. We introduce a string family where $\nu$ is asymptotically strictly smaller than both $\ell$ and $b$, which shows that NU-systems are indeed relevant and positions $\nu$ as the unique smallest reachable repetitiveness measure to date that captures both kinds of repetitiveness in non-trivial ways.

4. We study various ways of simplifying L-systems and show that, in most cases, we end up with a weaker repetitiveness measure. We also study some of those weaker variants of $\ell$, which can be of independent interest.

Overall, our results contribute to understanding how to measure repetitiveness and how to exploit it in order to build better compressors. We summarize the state of $\ell$ and $\nu$ after this work in Figure 1.

2 Basic concepts

In this section we explain the basic concepts needed to understand the rest of the paper, from strings and morphisms to relevant repetitiveness measures.

2.1 Strings

A (finite) string $w$ is a mapping from $\mathbb{Z}$ to $\Sigma$, and its length is called $\omega$, which is greater than $n$ for any $n \in \mathbb{Z}^+$. It is possible to define the concatenation $x \cdot y$ if $x$ is finite and $y$ infinite. The concepts of substring, prefix, and suffix carry over to infinite strings, with proper prefixes always being finite and suffixes always being infinite. The notations $w[i], w[i : j], w[: i]$ and $w[i :] = w[i]w[i+1] \ldots$ also carry over to infinite strings.
2.2 Morphisms

The set $\Sigma^*$ together with the (associative) concatenation operator and the (identity) string $\varepsilon$ form a monoid structure $(\Sigma^*, \cdot, \varepsilon)$. A morphism on strings is a function $\varphi : \Sigma^*_1 \rightarrow \Sigma^*_2$ satisfying $\varphi(x \cdot y) = \varphi(x) \cdot \varphi(y)$ for all $x, y$ (i.e., a function preserving the monoid structure), where $\Sigma_1$ and $\Sigma_2$ are alphabets. To define a morphism on strings, it is sufficient to define how it acts over the symbols in its domain. The pairs $(a, \varphi(a))$ for $a \in \Sigma_1$, usually denoted $a \rightarrow \varphi(a)$, are called the rules of the morphism, and there are $|\Sigma_1|$ of them. If $\Sigma_1 = \Sigma_2$, then the morphism is called an endomorphism.

Let $\varphi : \Sigma^*_1 \rightarrow \Sigma^*_2$ be a morphism on strings. Some useful definitions are $\text{width}(\varphi) = \max_{a \in \Sigma_1} |\varphi(a)|$ and $\text{size}(\varphi) = \sum_{a \in \Sigma_1} |\varphi(a)|$. A morphism is non-erasing if $\forall a \in \Sigma_1, |\varphi(a)| > 0$, expanding if $\forall a \in \Sigma_1, |\varphi(a)| > 1$, $k$-uniform if $\forall a \in \Sigma_1, |\varphi(a)| = k > 1$, and it is a coding if $\forall a \in \Sigma_1, |\varphi(a)| = 1$ (sometimes called a 1-uniform morphism).

Let $\varphi : \Sigma^* \rightarrow \Sigma^*$ be an endomorphism. Then $\varphi$ is prolongable on a symbol $a$ if $\varphi(a) = ax$ for some string $x \neq \varepsilon$. If this is the case, then for each $i, j$ with $0 \leq i \leq j$, it holds that $\varphi^i(a)$ is a prefix of $\varphi^j(a)$, and $x = \varphi^\omega(a) = ax\varphi(x)\varphi^2(x)\ldots$ is the unique infinite fixed-point of $\varphi$ starting with the symbol $a$. An infinite string $w = \varphi^\omega(a)$ that is the fixed-point of a morphism is called a purely morphic word, its image under a coding $x = \tau(w)$ is called a morphic word, and if the morphism $\varphi$ is $k$-uniform, then $x$ is said to be $k$-automatic.

2.3 Repetitiveness measures

A repetitiveness measure $\mu$ is a function that arguably captures the degree of repetitiveness of strings. Repetitiveness is an intuitive and elementary concept, yet it is still subject to debate. In general, a repetitive string is understood as one containing many copies of the same substrings. The more repetitive is a string $w$, the smaller the value $\mu(w)$ should be.

If we can represent every string $w[1 : n]$ within space $O(\mu(w))$ (where the asymptotics refer to $n$), then we say the measure $\mu$ is reachable. Space is usually measured in $\Theta(\log n)$-bit words following the conventions of the transdichotomous RAM model of computation. Hence, $O(\mu(w))$ space means $O(\mu(w) \log n)$ bits. We can represent any symbol in the alphabet of $w[1 : n]$ using a constant number of words as long as $|\Sigma| = O(n^d)$ for some $d \geq 0$.

A repetitiveness measure $u_1$ is smaller or lower-bounds another measure $u_2$ if $u_1(w) = O(u_2(w))$ for every $w[1 : n] \in \Sigma^*$. If, in addition, there is an infinite string family $F \subseteq \Sigma^*$ where $u_1(w) = o(u_2(w))$ for every $w[1 : n] \in F$, we say that $u_1$ is strictly smaller or strictly lower-bounds $u_2$. Two repetitiveness measures $u_1$ and $u_2$ are equivalent if each one is smaller than the other, and uncomparable if none is (i.e., $u_1 = o(u_2)$ on a string family $F_1$ and $u_2 = o(u_1)$ on another string family $F_2$).

In the following, we explain the most relevant repetitiveness measures to be considered in the rest of the paper. For a more in-depth review, see a recent survey [17].

Grammar-based measures

There exist several compressors, and measures of repetitiveness associated with their size, that build on context-free grammars.

A straight-line program (SLP) is a deterministic context-free grammar $G$ in Chomsky Normal Form whose language is a singleton $\{w\}$. We denote the string generated by the SLP as $exp(G) = w$, and extend this notation to the unique strings generated by the non-terminals of the grammar. The measure $g$ is defined as the size of the smallest SLP $G$ generating $w$. Finding the smallest SLP is an NP-complete problem [3], although there exist algorithms providing log-approximations [7, 24].
A measure based on context-free grammars that strictly lower-bounds $g$ is $g_{rt}$, the size of the smallest run-length SLP (RLSLP) generating $w$ [20]. RLSLPs allow constant-size rules of the form $A \rightarrow B^n$ for $n > 1$, which can make a noticeable difference in some string families like $\{a^n | n \geq 0\}$, where $g = \Theta(\log n)$, but $g_{rt} = O(1)$.

A collage system is an RLSP that, in addition, supports rules of the form $A \rightarrow B[i:j]$ for some $i, j \in [1, \exp(B)]$. These mean that $\exp(A) = \exp(B)[i:j]$. The size $c$ of the smallest collage system [10] strictly lower-bounds $g_{rt}$.

Parsing-based measures

A parsing of size $k$ produces a factorization of a string $w$ into non-empty phrases, i.e., $w = w_1w_2...w_k$ where $w_i \in \Sigma^+$ for $i \in [1,k]$. Several compressors work by parsing $w$ in a way that storing summary information about the phrases enables recovering $w$.

The Lempel-Ziv (LZ) parsing processes a string greedily from left to right, always forming the longest phrase that has a copy (called a source) starting inside some previous phrase or forming an explicit phrase of length 1 otherwise [14]. The source can overlap the new phrase. The LZ-no parsing, instead, does not allow the source overlap the new phrase. The LZ-end parsing [13] requires, in addition, that the source ends at a previous phrase boundary. All of these parsings can be constructed in linear time, and their number of phrases are denoted by $z$, $z_{na}$, and $z_e$, respectively. While $z$ and $z_{na}$ are optimal among the parsings satisfying their respective conditions, this is not always the case for $z_e$. The optimal factorization where each phrase $w_{i+1}$ appears as a suffix of $w_1...w_j$ for some $j \leq i$ is denoted by $z_{end}$. Because of the optimality of $z$, $z_{na}$, and $z_{end}$, it holds that $z \leq z_{na} \leq z_{end} \leq z_e$ for every string.

A bidirectional macro scheme (BMS) [26] is any parsing where each phrase of length greater than 1 has a copy starting at a different position in such a way that the original string can be recovered following these pointers (assuming that the phrases of length 1 store their symbol explicitly). The measure $b(w)$ is defined as the size of the smallest BMS for $w$. It strictly lower-bounds all the other reachable repetitiveness measures [18], except for the ones we focus on in this paper, $\ell$ and $\nu$ [19]. On the other hand, $b$ is NP-hard to compute [5].

Another interesting parsing-based measure is the size of the greedy lexicographic parsing of $w$, denoted as $\nu(w)$ [18]. This parsing processes $w$ from left to right, taking as the next phrase the longest common prefix between the unprocessed part of the string and a lexicographically smaller suffix of the processed part (a unique symbol $\$, smaller than the others, is assumed to exist at the end of $w$). It forms an explicit phrase of length one if the longest common prefix is empty or no predecessor exists.

Burrows-Wheeler transform

The Burrows-Wheeler transform (BWT) [2] is a reversible transformation that usually makes a string more compressible. It is obtained by concatenating the last symbols of the sorted suffixes of $w\$, where $\$ is a symbol smaller than any symbol appearing in $w$. The BWT tends to produce long runs of the same symbol when a string is repetitive, and these (maximal) runs can be compressed into one symbol in the alphabet $\{(a,k) | a \in \Sigma, k \in [1,n]\}$ using run-length encoding (rle). A repetitiveness measure based on this idea is defined as $r(w) = |rle(BW(T(w))|$. Although $r$ is not ideal as a repetitiveness measure [6], its size can be bound in terms of $z$ [8]. It has many practical applications representing repetitive sequences in Bioinformatics because of its indexing power [4].
L-Systems for Measuring Repetitiveness

String attractors
Kempa and Prezza [9] introduced the notion of string attractor as a unifying framework and lower bound for dictionary-based compressors.

Let $w$ be a string of length $n$. A string attractor for $w$ is a set of positions $\Gamma \subseteq [1, n]$ such that for each substring $w[i : j]$ of $w$, there exist integers $i', j' \in [1, n]$ and $k \in \Gamma$, such that $w[i : j] = w[i' : j']$ and $i' \leq k \leq j'$. That is, every substring of $w$ has a copy covering a position in $\Gamma$. The measure $\gamma(w)$ is defined as the size of the smallest string attractor for $w$.

Computing $\gamma$ is an NP-complete problem. The measure $\gamma$ is a lower bound to $b$ and is also strictly smaller than $b$ when considering the infinite family of Thue-Morse words [1]. On the other hand, it is unknown if $\gamma$ space or even $o(\gamma \log(n/\gamma))$ space is reachable.

Substring complexity
Let $F_w(k)$ be the set of distinct substrings of $w$ of length $k$. The complexity function of $w$ is defined as $P_w(k) = |F_w(k)|$. Kociumaka et al. [12] introduced a repetitiveness measure based on the complexity function, defined as $\delta(w) = \max \{P_w(k)/k | k \in [1..|w|]\}$.

The measure $\delta$ has several nice properties: it is computable in linear time, monotone, insensitive to reversals, resistant to small edits on $w$, can be used to construct $O(\delta \log(n/\delta))$-space representations supporting efficient access and pattern matching queries [12, 11], and is a lower bound to almost every other theoretical or ad-hoc repetitiveness measure considered in the literature, including $\gamma$. On the other hand, $o(\delta \log(n/\delta))$ space is unreachable [12].

3 The measure $\ell$

The class of CPD0L-systems is a variant of the original L-systems, the parallel grammars without terminals defined by Aristid Lindenmayer to model cell divisions in the growth of plants and algae [15, 16].

Formally, a CPD0L-system is a 4-tuple $L = (\Sigma, \varphi, \tau, s)$, where $\Sigma$ is the alphabet, $\varphi$ is the set of rules (a non-erasing endomorphism on $\Sigma^*$), $\tau$ is a coding on $\Sigma^*$, and $s \in \Sigma^*$ is the axiom. The system generates the sequence $(\tau(\varphi^d(s)))_{d \in \mathbb{N}}$. The “D0L” stands for deterministic L-system with 0 interactions. The “P” stands for propagating, which means that it has no $\varepsilon$-rules. The “C” stands for coding, which means that the system is extended with a coding. To define a compressor based on CPD0L-system, we extend them to 6-tuples by fixing $d$ and using another parameter $n$, so we can uniquely determine a string of the sequence generated by a system and then extract a prefix from it. For simplicity, in the rest of this paper, we refer to these extended CPD0L-systems just as L-systems.

**Definition 1 (L-systems).** An L-system is a 6-tuple $L = (\Sigma, \varphi, \tau, s, d, n)$ where $\Sigma$ is the alphabet, $\varphi$ is the set of rules (an endomorphism on $\Sigma^*$), $\tau$ is a coding on $\Sigma^*$, $s \in \Sigma^*$ is the axiom, and $d$ and $n$ are two non-negative integers. The string generated by $L$ is $w = \tau(\varphi^d(s))[1 : n]$.

We now define the size of an L-system and the measure $\ell$.

**Definition 2 (Measure $\ell$).** The size of an L-system $L = (\Sigma, \varphi, \tau, s, d, n)$ is $\text{size}(L) = \text{size}(\varphi) + |s| + |\Sigma| + 2$. The measure $\ell(w)$ is defined as the size of the smallest L-system generating $w$.

The size of an L-system accounts for the lengths of the right-hand sides of its rules, the length of the axiom, the coding $\tau$, and the values $d$ and $n$, so we can effectively represent the system using $O(\text{size}(L))$ space. Hence, the measure $\ell$ is reachable.
As a convention, we always assume that \( d = n^{O(1)} \) and that \( \Sigma = n^{O(1)} \). Otherwise, we could need too many words to represent the integer \( d \) or the symbols of the alphabet.

A finer-grained analysis about the number of bits needed to represent an L-system of size \( \ell \) yields \( O(\ell \log |\Sigma| + \log n) \) bits, the second term corresponding to \( d \) and \( n \); note that \( \Sigma \) contains the alphabet of \( w \).

An important result about \( \ell \) is that it always holds that \( \ell = O(g) \) [19]. More importantly, sometimes \( \ell = o(\delta) \), which implies that \( \delta \) is not a lower bound for \( \ell \), and questions \( \delta \) as a definitive measure of repetitiveness.

To understand the particularities of \( \ell \), we study several classes of L-systems with different restrictions and define measures based on them. We define the measure \( \ell_e \) that restricts the morphism to be expanding. The measure \( \ell_u \) restricts the morphism to be \( k \)-uniform for some \( k > 1 \). The variant \( \ell_m \) forces the morphism of the system to be \( a \)-prolongable for some symbol \( a \) and the axiom to be \( s = a \). The variant \( \ell_d \) essentially removes the coding by forcing it to be the identity function. Finally, \( \ell_p \) refers to the intersection of \( \ell_m \) and \( \ell_d \), and \( \ell_a \) refers to the intersection of \( \ell_m \) and \( \ell_u \), which intuitively perform well in prefixes of purely morphic words and prefixes of automatic sequences, respectively.

\[ \text{Definition 3.} \quad \text{An L-system} \ (\Sigma, \varphi, \tau, s) \ \text{is a-prolongable if there exists a symbol} \ a \ \text{such that} \ s = a \ \text{and} \ a \rightarrow ax \ \text{with} \ x \neq \varepsilon. \ \text{An L-system is prolongable if it is a-prolongable for some symbol} \ a. \]

\[ \text{Definition 4 (}\ell\text{-variants).} \ \text{We define the following} \ell\text{-variants} \]
1. The variant \( \ell_e \) denotes the size of the smallest L-system generating \( w \), satisfying that all its rules have a size at least 2.
2. The variant \( \ell_m \) denotes the size of the smallest prolongable L-system generating \( w \).
3. The variant \( \ell_d \) denotes the size of the smallest L-system generating \( w \), satisfying that \( \tau \) is the identity function.
4. The variant \( \ell_u \) denotes the size of the smallest L-system generating \( w \), satisfying that all its rules have the same size, at least 2.
5. The variant \( \ell_p \) denotes the size of the smallest prolongable L-system generating \( w \), satisfying that \( \tau \) is the identity function.
6. The variant \( \ell_a \) denotes the size of the smallest prolongable L-system generating \( w \), satisfying that all its rules have the same size, at least 2.

It is known that different classes of L-systems produce different classes of languages [21]. Some of these classes also differ in the factor complexity of the sequences they can generate [22].

It is interesting to understand how these differences in terms of expressive power and factor complexity translate into compression power.

We defer the study of \( \ell \)-variants to Section 7. We define them early, as some of our results relating \( \ell \) to better-established measures in Sections 4 and 5 also apply for some of the \( \ell \)-variants.

### 4 Breaking the \( \delta \)-lower-bound for repetitiveness

It is known that the repetitiveness measure \( \delta \) is a (strict) lower bound to all the other repetitiveness measures [17]. It is also known that \( \delta \) is a lower bound to \( k \)-th order empirical entropy, which plays a role in several compressors [17]. This makes \( \delta \) an asymptotic lower bound to the size of almost every existing compressor and compressibility measure.
Certainly, we cannot expect to find a reachable measure upper-bounded by $O(\delta)$ because $\delta$-space is unreachable, as shown by Kociumaka et al. [12]. Still, it could be possible to design measures capturing repetitiveness and going below $\delta$ in some restricted but relevant scenarios. In this context, we raise the following question:

Can we find a competitive and reachable repetitiveness measure achieving space lower than $\delta$ on some restricted but relevant cases?

It is not difficult to design a trivial measure breaking the $\delta$-lower-bound for some specific string families. We also require, however, this measure to make sense, be reachable, and be competitive, the latter meaning that it is at least as good as $z$, $g$, or $r$ (i.e., the most popular reachable measures in practice) in terms of space.

The repetitiveness measure $\ell$ was designed with the conditions above in mind. As we already mentioned, $\ell$ cannot lower-bound $\delta$ because $\ell$ is a reachable measure.

Lemma 5 ([19, Theorem 4]). There exist string families where $\ell = \Omega(\delta \log n)$.

It was shown that $\ell$ is a competitive repetitiveness measure: the smallest L-system for a string is always asymptotically smaller than the smallest grammar (their proof applies to the variant $\ell_d$ as well). This shows that the measure $\ell$ is always reasonable for repetitive strings:

Lemma 6 ([19, Theorem 6]). It always holds that $\ell = O(g)$.

On the other hand, they [19] showed a string family satisfying that $\delta = \Omega(\ell \log n)$, and conjectured that this gap was the maximum possible, that is, that the lower bound $\ell = \Omega(\delta / \log n)$ holds for any string family. We now disprove this conjecture. We show a string family where $\delta$ is $\Theta(\sqrt{n})$ times bigger than the size $\ell$ of the smallest L-system.

Lemma 7. There exists a string family where $\delta = \Theta(\ell \sqrt{n})$.

Proof. Consider a $c$-prolongable L-system $L_d = (\Sigma, \varphi, \tau, s, d, n)$, where

$$
\begin{align*}
\Sigma &= \{a, b, c\} \\
\varphi &= \{a \to a, b \to ab, c \to cb\} \\
\tau &= \{a \to a, b \to b, c \to c\} \\
s &= c \\
n &= 1 + \frac{(d-1)d}{2} + d
\end{align*}
$$

for any $d \geq 0$. By iterating the morphism $\varphi$ we obtain the words

$$
\begin{align*}
\varphi^0(c) &= c \\
\varphi^1(c) &= cb \\
\varphi^2(c) &= cbab \\
\varphi^3(c) &= cbabaab \\
\varphi^4(c) &= cbabaabaaaaab \\
\varphi^5(c) &= cbabaabaaaaab
\end{align*}
$$

and so on, from which we extract as a prefix the whole word (depending on the value of $d$ chosen). It is easy to check by induction that for each $d \geq 0$, the string generated by the system $L_d$ is $s_d = c \prod_{i=0}^{d-1} a^i b$ and it has length $1 + \frac{(d-1)d}{2} + d$. 


$s_3 = \text{c b a b a a b}$

$s_6 = \text{c b a b a a b a a a b a a a a b a a a a a b}$

**Figure 2** All the substrings of length 6 of the string $s_6$ of Lemma 7 starting inside some position $i \leq |s_3| = 7$ are distinct, because the runs of a’s considered have all different and increasing lengths, and $d$ is big enough. The last of the substrings considered is underlined. Extending these substrings one position to the left yields $|s_3|$ different strings of length 7, so the claim holds for even and odd values of $d \geq 2$.

It holds that $\ell$ is $\Theta(1)$ in this family. The system is essentially the same for every string in the family. The only changes are the integers $d$ and $n$, which always fit in constant space.

On the other hand, the first $|s_{|d|/2}| = 1 + (\lfloor |d|/2 \rfloor - 1)(\lfloor |d|/2 \rfloor )/2 + |d|/2$ substrings of length $d$ of $s_d$ (for $d \geq 2$) are completely determined by the b’s they cross, and the number of a’s at their extremes, so they are all distinct. An example can be seen in Figure 2.

This gives the lower bound $\delta = \Omega(d) = \Omega(\sqrt{n})$. The upper bound $O(\sqrt{n})$ holds trivially for run-length grammars as the strings considered have $\Theta(\sqrt{n})$ runs of a’s followed by b’s, so $\delta = \Theta(\sqrt{n})$. Thus $\delta = \Theta(\ell \sqrt{n})$ in this string family.

This string of Lemma 7 is easy to describe yet hard to represent with copy-paste mechanisms. Intuitively, the simplicity of the sequence relies on the fact that many substrings can be described in terms of previous ones, so it is arguably highly repetitive, though not via copy-paste. The repetitiveness in this family is better captured by an L-system, instead.

## 5 Uncomparability of $\ell$ with other repetitiveness measures

As a corollary of Lemmas 5 and 7 (and also mentioned in previous work [19]), we obtain that $\ell$ and $\delta$ are uncomparable as repetitiveness measures.

**Corollary 8.** The measures $\ell$ and $\delta$ are uncomparable.

Moreover, this is also true for the variant $\ell_p$ because, in Lemma 7, we considered a prolongable L-system with the identity function as the coding. As we prove later in Section 7, the variant $\ell_p$ is, in general, far from ideal for measuring repetitiveness, so the fact that $\delta$ is uncomparable to this weak variant is even more surprising.

A natural question is then to identify which other measures are also uncomparable to $\ell$ (and $\ell_p$). We show in this section that this holds for almost any other repetitiveness measure. To do so, we first recall the string family defined by Kociumaka et al. [12], satisfying that it needs $\Omega(\log^2 n)$ bits to be represented with any method. This string family will be crucial in the following proofs.

**Definition 9 (12).** The string family $\mathcal{K}$ is formed by all the infinite strings $s$ over $\{a, b\}$ constructed as follows:

2. For any $i \geq 2$, choose a position $j_i$ in $[2 \cdot 4^{i-2} + 1, 4^{i-1}]$. Then, $s[j_i] = b$.
3. If $j > 1$ and $j \neq j_i$ for any $i \geq 2$, then $s[j] = a$.

The family $\mathcal{K}_n$ for $n \geq 0$ is formed by all the prefixes of length $n$ of some string in $\mathcal{K}$.

It is easy to see that the strings in the family $\mathcal{K}_n$ have $\Theta(\log n)$ symbol b’s. Also, note that with the possible exception of the first two positions, there are no consecutive b’s.
Now we are ready to prove that, in general, it does not hold that $\ell = O(g_{rl})$, making L-systems uncomparable to RLSLPs.

Lemma 10. There exists a string family where $\ell = \Omega(g_{rl} \log n / \log \log n)$.

Proof. Consider the string family $K_n$ needing $\Omega(\log^2 n)$ bits (or $\Omega(\log n)$ space) to be represented with any method [12]. Strings in $K_n$ have $O(\log n)$ runs of $a$’s separated by $b$’s, so it is easy to see that $g_{rl} = O(\log n)$ in this family. Because of this, and because $g_{rl}$ is a reachable measure, it holds that $g_{rl} = \Theta(\log n)$ in $K_n$. On the other hand, the minimal L-system for a string in this family can be represented with $O(\ell \log |\Sigma| + \log n) \subseteq O(\ell \log \ell + \log n)$ bits, which must be in $\Omega(\log^2 n)$ bits because the L-system is also reachable. It follows that $\ell = \Omega(\log^2 n / \log \log n)$; otherwise,

$$\ell \log \ell = o((\log^2 n / \log \log n) \log(\log^2 n / \log \log n))$$

$$= o(\log^2 n),$$

which contradicts $\ell$ being reachable. Thus, $\ell = \Omega(g_{rl} \log n / \log \log n)$ in this string family. ▶

The same result holds for LZ-like parsings. Even the greedy LZ-End parsing (the largest of them) can be asymptotically smaller than $\ell$ in some string families.

Lemma 11. There exists a string family where $\ell = \Omega(z_e \log n / \log \log n)$.

Proof. Take each string in $K_n$ and prepend $a^n$ to it. This new family of strings still needs $\Omega(\log^2 n)$ bits to be represented with any method because the size of the family is the same, and $n$ just doubled. Just as in Lemma 10, it holds that $\ell = \Omega(\log^2 n / \log \log n)$ in this family. On the other hand, the LZ-End parsing needs $\Theta(\log n)$ phrases only to represent the prefix $a^n b$, and then for each run of $a$’s followed by $b$, its source is aligned with $a^n b$, so $z_e = \Theta(\log n)$. Thus, $\ell = \Omega(z_e \log n / \log \log n)$. ▶

The same result also holds for the number of equal-letter runs of the Burrows-Wheeler transform of a string.

Lemma 12. There exists a string family where $\ell = \Omega(r \log n / \log \log n)$.

Proof. Consider the family $K_n$ again. Clearly $r = \Omega(\log n)$, because $r$ is reachable. Because a string in this family has $O(\log n)$ $b$’s, its BWT has also $O(\log n)$ runs of $a$’s separated by $b$’s (or the unique $\$). Thus, $r = \Theta(\log n)$ and $\ell = \Omega(r \log n / \log \log n)$ in this string family. ▶

We conclude that the measure $\ell$ is uncomparable to almost every other repetitiveness measure. We summarize these results in the following theorem.

Theorem 13. The measure $\ell$ is incomparable to the repetitiveness measures $\delta$, $b$, $v$, $c$, $g_{rl}$, $z$, $z_{no}$, $z_{end}$, $z_e$, and $r$.

Proof. There exist string families where $\ell = o(\delta)$. In these families, it holds $\ell = o(\mu)$ where $\mu$ is any of the measures considered above, because $\delta$ is a lower bound to all of them. On the other hand, all the measures above are upper-bounded by at least one of $z_e$, $g_{rl}$, or $r$, which by Lemmas 10, 11, and 12, respectively, can be asymptotically smaller than $\ell$ for some string families. ▶

This shows that $\ell$, although reachable and competitive as a repetitiveness measure, captures the regularities in strings in a form that is largely orthogonal to other repetitiveness measures. As the underlying regularities captured by $\ell$ and the other measures are apparently different, we try to combine them to obtain more powerful measures/compressors.
6 NU-systems and the measure $\nu$

A NU-system [19] is a tuple $N = (V, R, \Gamma, s, d, n)$ that generates a unique string in a similar way to an L-system. The key difference is that on the right-hand side of its rules, a NU-system is permitted to have special symbols of the form $a(k)[i : j]$, whose meaning is to generate the $k$-th level from $a$, then extract the substring starting at position $i$ and ending at position $j$, and finally apply the coding to the resulting substring.

The indices in a NU-system (e.g., levels, intervals) must be less or equal to $n$ to fit in a $\Theta(\log n)$-bits word. Also, the NU-system must not produce any loops when extracting a prefix from some level, which is decidable to detect. The size of a NU-system is defined analogously to the size of L-systems, with the extraction symbols $a(k)[i : j]$ being symbols of length 4. The measure $\nu$ is defined as the size of the smallest NU-system generating $w$.

It holds that $\nu = O(\ell)$ and $\nu = O(b)$ [19]. Moreover, there exist families where both asymptotic bounds are strict. We now show that NU-systems exploit the features of L-systems and macro schemes in a way that, for some string families, can reach sizes that are unreachable for both L-systems and macro schemes independently.

\textbf{Theorem 14.} There exists a family of strings where $\nu = o(\min(\ell, b))$.

\textbf{Proof.} Let $\mathcal{K}_m$ be the family of strings defined by Kociumaka et al. [12], needing $\Omega(\log^2 m)$ bits to be represented with any method, but over the alphabet $\{0, 1\}$. We construct a new family $\mathcal{F} = \{x \cdot y[m] | x \in \mathcal{K}_m\}$, where $y$ is the infinite fixed point generated by the c-prolongable L-system with the identity function as the coding, utilized in Lemma 7.

Let $n = 2m$. As shown in Lemma 10, it holds that $\ell = \Omega(\log^2 n / \log \log n)$ in this family. On the other hand, $b = \Omega(\sqrt{n})$, because $d = \Omega(\sqrt{n})$ on prefixes of $y$, and the alphabets between the prefix in $\mathcal{K}_m$ and $y[m]$ are disjoint.

Let $x$ be a string in $\mathcal{K}_m$ with $k$ symbols 1. Let $i_j$ be the number of 0’s in $x$ between the $(j - 1)$-th 1 and the $j$-th 1, for $j \in [2, k]$. Also, let $i_1$ and $i_{k+1}$ be the number of 0’s at the left and right extremes of $x$. We construct a NU-system $N = (V, R, \Gamma, s, d, n)$ as follows:

$$V = \{0, 1, a, b, c\}$$
$$R = \{0 \rightarrow 00, 1 \rightarrow 1, a \rightarrow a, b \rightarrow ab, c \rightarrow cb\}$$
$$\Gamma = \{0 \rightarrow 0, 1 \rightarrow 1, a \rightarrow a, b \rightarrow b, c \rightarrow c\}$$
$$s = 0(m)[i_1]10(m)[i_2]0(m)[i_3]10(m)[i_4]c(m)[m]$$
$$d = 0$$
$$n = 2m$$

By construction, this NU-system generates the string $x \cdot y[m]$ of length $n$, and its axiom has size $4(k + 2) + k$, where $k = \Theta(\log n)$. Hence, it holds that $\nu = O(\log n)$ for these strings. Thus, $\nu = o(\min(\ell, b))$ in the family $\mathcal{F}$ we constructed. \hfill $\blacktriangle$

NU-systems can then be smaller representations than those produced by any other compression method exploiting repetitiveness. This shows that combining copy-paste mechanisms with iterated morphisms is an effective way of improving compression from a theoretical point of view.

On the other hand, though computable, no efficient decompression scheme has been devised for NU-systems. In turn, finding the smallest NU-system is very likely an NP-complete problem, and probably very difficult to even approximate.
L-systems are weaker than $\ell$

In this section we show that the features we include in the L-systems used in the definition of $\ell$ are necessary to obtain a competitive repetitiveness measure; removing any of them yields an inherent loss in compression power.

We start by showing that $\ell$ can be asymptotically strictly smaller than $\ell_m$. That is, restricting L-systems to be prolongable yields a weaker measure.

**Lemma 15.** There exists a string family where $\ell = o(\ell_m)$.

**Proof.** Let $F = \{a^{n-1}b \mid n \geq 1\}$. Clearly, $\ell$ is constant in this string family: the L-system $L_n = (V, \varphi, \tau, s, d, n)$ where $V = \{a, b\}$, $\varphi = \{a \to a, b \to ab\}$, $\tau = \{a \to a, b \to b\}$, $s = b$, and $d = n - 1$ produces each string in $F$ by changing only the value of $n$ and $d$ accordingly. Note that these L-systems are not prolongable on the axiom.

For the sake of contradiction, suppose that $\ell_m = O(1)$ in $F$. Let $L_n = (\Sigma_n, \varphi_n, \tau_n, c, d_n, n)$ be the the smallest $c$-prolongable system generating $a^{n-1}b$. Because $\ell_m = O(1)$, there exists a constant $C$ satisfying that $|\Sigma_n| < C$ and $\text{width}(\varphi_n) < C$ for every $n$. Observe that it is only necessary to have one symbol $c' \in \Sigma_n$ with $\tau_n(c') = b$ because there is only one $b$ in $a^{n-1}b$, so w.l.o.g. assume that $b \in \Sigma_n$ and $\tau_n(b) = b$. As the system is $c$-prolongable, each level is a prefix of the next one. This implies that the morphism should be iterated until $b$ appears for the first time, and then we can safely extract the prefix. This must happen in the first $C$ iterations of the morphism; otherwise, $b$ is not reachable from $C$ (i.e., if an iteration does not yield a new symbol, then no new symbols will appear since then, and there are less than $C$ symbols). But in the first $C$ iterations, we cannot produce a string longer than the constant $C^C$. For sufficiently large $n$, this implies that the symbol $b$, if it is reachable, will appear for the first time before the $n$-th position, which is a contradiction. ▶

Because we used the identity coding in the proof above, we can obtain the following corollary.

**Corollary 16.** There exists a string family where $\ell_d = o(\ell_m)$.

We can prove a similar result for uniform morphisms.

**Proposition 17.** There exists a string family where $\ell_u = o(\ell_m)$.

**Proof.** It is not difficult to see that $\ell_u$ is constant in the family $\{a^{2^k}b \mid k \geq 0\}$: consider the axiom $s = ab$ and the rules $a \to aa$, $b \to bb$, the level $d = k$ and the prefix length $n = 2^k + 1$. A similar argument to the one of Lemma 15 yields that $\ell_u = o(\ell_m)$ for this other string family. ▶

Further, we can find a concrete asymptotic gap between $\ell$ and $\ell_m$ in the string family of the proof of the previous lemma.

**Lemma 18.** There exists a string family where $\ell_m = \Omega(\ell \log n / \log \log n)$.

**Proof.** Let $F = \{a^{n-1}b \mid n \geq 1\}$. Recall that $\ell = O(1)$ in this family. Let $k = |\Sigma|$ and $t = \text{width}(\varphi)$ obtained from the morphism of the smallest $c$-prolongable system generating $a^{n-1}b$ (we assume again that the only symbol mapped to $b$ by the coding is $b$). In the first $k$ iterations, $b$ must appear (as in the previous proof) and cannot be deleted in the following levels, so it cannot appear before position $n$. Hence, $tk \geq n$, which implies
\( k \geq \log n \). By definition, \( \ell_m \geq k \geq \log_t n \) and \( \ell_m \geq t \), so \( \ell_m \geq \max(t, \log_t n) \). The solution to the equation \( t = \log_t n \) is the smallest value that \( \max(t, \log_t n) \) can take for \( t \in [2..n] \). This value is \( \Omega(\log(n/W(\log n))) \) where \( W(x) \) is the Lambert W function, and it holds that \( W(\log n) = \Theta(\log \log n) \). Therefore, \( \ell_m = \Omega(\ell \log n / \log \log n) \) in this string family. ▶

As a corollary, we obtain the following result.

**Corollary 19.** There exists a string family where \( \ell_m = \Omega(\ell_d \log n / \log \log n) \).

We now show that if we remove the coding from prolongable L-systems, which corresponds to the variant \( \ell_p \), we end with a much worse measure. We change the usual alphabet for clarity of presentation.

**Lemma 20.** There exists a string family where \( \ell_p = \Omega(\ell_m \sqrt{n}) \).

**Proof.** We prove that \( \ell_p = \Theta(n) \) whereas \( \ell_m = O(\sqrt{n}) \) on \( F = \{0^{n-1}1 \mid n \geq 2 \} \). Any prolongable morphism with an identity coding generating \( 0^{n-1}1 \) must have the rule \( 0 \rightarrow 0^n1 \), which implies \( \ell_p = \Theta(n) \). The reason is that if the system is prolongable, but it has no coding, then the axiom must be \( 0 \), and in the prolongable rule \( 0 \rightarrow 0w \), if \( |\varphi(0)| \leq n \), then the non-empty string \( w \) could only contain \( 0 \)'s and \( 1 \)'s, otherwise undesired symbols would appear in the final string because the starting level is a prefix of the final level. If \( w \) does not contain \( 1 \)'s, then \( 1 \) is unreachable from \( 0 \). If \( w \) contains a \( 1 \), then the first of them should be at position \( n \).

On the other hand, we can construct an \( a \)-prolongable morphism, with \( \tau(1) = 1 \) and \( \tau(a) = 0 \) for every other symbol \( a \neq 1 \) as follows: Let \( n - 1 = k[\sqrt{n - 1}] + j \) with \( \lceil \sqrt{n - 1} \rceil > 3, k > 1, 0 \leq j < \lceil \sqrt{n - 1} \rceil \) ( \( k \) and \( j \) integers). We can assume \( n \) is sufficiently big so the requirements are satisfied. Then, define the following rules

\[
\begin{align*}
&\quad a \rightarrow ab \\
&\quad b \rightarrow a^{k-1}d \\
&\quad c \rightarrow 0[\sqrt{n-1}]^{-1} \\
&\quad d \rightarrow 0[\sqrt{n-1}]^{-3+j}1.
\end{align*}
\]

The first four levels are

\[
\begin{align*}
\varphi^0(a) &= a \\
\varphi^1(a) &= ab \\
\varphi^2(a) &= abc^{k-1}d \\
\varphi^3(a) &= abc^{k-1}d0((\sqrt{n-1})^{-1}(k-1)0[\sqrt{n-1}]^{-3+j}1,
\end{align*}
\]

and it holds that

\[
|\varphi^3(a)| = 3 + (k - 1) + (\lceil \sqrt{n - 1} \rceil - 1)(k - 1) + (\lceil \sqrt{n - 1} \rceil - 3 + j) + 1 = n.
\]

Hence, \( \tau(\varphi^3(a)) = 0^{n-1}1 \). The system \( L = \{a, b, c, d, 0, 1\}, \varphi, \tau, a, 3, n\} \) generates \( 0^{n-1}1 \) as required for \( n \) bigger than some constant. The size of the system is clearly \( O(\sqrt{n}) \). Thus, the claim holds. ▶

By using the same family above, the following corollary holds.

**Corollary 21.** There exists a string family where \( \ell_p = \Omega(\ell_d \sqrt{n}) \).
It is surprising that this weak measure \( \ell_p \) can be much smaller than \( \delta \) for some string families. This can be deduced from Lemma 7. On the other hand, it does not hold that \( \ell_p = O(g) \) for any string family, because \( g = \Theta(\log n) \) on \( \{0^{n-1}1 \mid n \geq 1\} \).

**Corollary 22.** The variant \( \ell_p \) is incomparable to the measures \( \delta \) and \( g \).

If we restrict L-systems to be expanding, that is, with all its rules having a length of at least 2, we also end with a weaker measure. This shows that, in general, it is not possible to transform L-systems into expanding ones without incurring an increase in size.

**Lemma 23.** There exists a string family where \( \ell = o(\ell_c) \).

**Proof.** Let \( \mathcal{F} = \{a^kba^{2k} \mid k \geq 0\} \). Clearly \( \ell \) is constant in \( \mathcal{F} \): the L-system \( \{\{a, b, c\}, \{a \rightarrow aa, b \rightarrow cb, c \rightarrow c\}, \{a \rightarrow a, b \rightarrow b, c \rightarrow a\}, ba, d = k, n = 2^k + k + 1\} \) produces \( a^kba^{2k} \) and stays constant-size as \( k \) (and \( d \) and \( n \)) grows.

Suppose that \( \ell_c \) is also constant in \( \mathcal{F} \). Then there is a constant \( C \) such that the minimal expanding L-systems generating the strings in this family have at most \( C \) rules, each one of length at most \( C \). Without loss of generality, assume that for each of these systems, the only symbol mapped to \( b \) by the coding is \( b \). Also, assume that the axiom is a single symbol \( a_0 \). Note that because the systems are expanding with rules of size at most \( C \), their level must be \( d \geq \log_C 2^k = \frac{k}{\log 2} \).

Let \( a_0, a_1, \ldots, a_d \) be the sequence of the first symbols of \( \varphi^i(a_0) \) for \( i \leq d \). By the pigeonhole principle, for sufficiently big values of \( k \) (and consequently big values of \( d \)), this sequence has a period of length \( q \) starting from \( a_p \), with \( p + q \leq C \leq d \). Then there exist indexes \( t \) and \( j \) such that \( t = d - jq \) and \( p \leq t < p + q \). By the \( q \)-periodicity of the sequence starting at \( a_t \), it is clear that \( \varphi^q(a_t) = a_tw \) for some \( w \neq \varepsilon \) (because the morphism is expanding), so \( \varphi^q \) is prolongable on \( a_t \). This implies that \( \varphi^q(a_t) \) is a prefix of \( \varphi^w(a_t) \) for \( i \leq d \). As before, if \( \mathcal{F} \) is reachable from \( a_t \) via \( \varphi^q \), that must happen in the first \( C \) iterations, so \( \varphi^C(a_t) \) contains a \( b \), and so does \( \varphi^q(a_t) \), which is a prefix of \( \varphi^w(a_t) \). This implies that \( \varphi^q(a_0) \) contains a \( b \) before position \( C^q, \) which is bounded by \( C^q, \) a contradiction for sufficiently long strings in the family. So it has to be that \( \mathcal{F} \) is not reachable via \( \varphi^q \) from \( a_t \), but this is also a contradiction for sufficiently long strings because \( \varphi^q(a_t) \) is a prefix of \( \varphi^q(a_0) \) of length at least \( 2^{d-1} = \omega(k) \), yielding too many symbols not mapped to \( b \) before the first \( b \) at level \( d \). Thus, \( \ell_c \) cannot be \( O(1) \) in \( \mathcal{F} \).

We summarize the results of this section in Figure 3. Overall, we have shown that imposing restrictions on the length of the rules of an L-system or forcing them to be prolongable wildly impacts their compression power. We have not yet found an example where \( \ell_d \) could be asymptotically smaller than \( \ell \), which would prove that the coding contributes to the measure \( \ell \) in a fundamental way (the purpose of the coding is to make \( \ell \) constant in the case of prefixes of general morphic words, but it is unknown if it is really needed). We conjecture that such a family exists and the coding is necessary.

## 8 Conclusions and open questions

The measure \( \ell \) is arguably a strong reachable repetitiveness measure, which can break the limits of \( \delta \) (a measure considered a stable lower bound for repetitiveness) by a wide margin (a factor of \( \sqrt{n} \)). On the other hand, however, \( \ell \) can be asymptotically weaker than the space reached by several compressors based on run-length context-free grammars, many Lempel-Ziv variants, and the Burrows-Wheeler transform. Only the size of context-free grammars is an upper bound to \( \ell \). This suggests that the self-similarity exploited by L-systems is mostly
Figure 3 Asymptotic relations between $\ell$-variants and other relevant measures. A black arrow from $v_1$ to $v_2$ means that it always holds that $v_1 = O(v_2)$. A double black arrow from $v_1$ to $v_2$ means that it also exists a string family where $v_1 = o(v_2)$. A dashed arrow from $v_1$ to $v_2$ means that there exists a family where $v_1 = o(v_2)$.

Independent of the source of repetitiveness exploited by other compressors and measures, which build on copy-paste mechanisms. We also show that several attempts to simplify or restrict L-systems lead to weaker measures.

A relevant question about L-systems is whether they can be useful for building compressed sequence representations that support direct access. More formally, can we build an $O(\ell)$-space representation of a string $w[1:n]$ providing random access to any position of the string in $O(\text{polylog } n)$ time? The closest result (as far as we know) is an algorithm designed by Shallit and Swart [25], which computes $\phi^d(a)[i]$ in time bounded by a polynomial in $|\Sigma|$, $\text{width}(\phi)$, $\log d$ and $\log i$. It uses more space and takes more time than our aim. The main bottleneck is having to store the incidence matrix of the morphism and compute its powers. As suggested by Shallit and Swart, this could be solved by finding closed forms for the growth functions (recurrences) of each symbol. If this approach were taken, these formulas should be easily described within $O(\ell)$ space.

In terms of improving compression, on the other hand, the recent measure $\nu$ [19] aims to unify the repetitiveness induced by self-similarity and by explicit copies. This measure is the smallest size of a NU-system, a natural way to combine L-systems (with minimum size $\ell$) with macro schemes (with minimum size $b \geq \delta$). In line with our finding that $\ell$ and $\delta$ are mostly orthogonal, we prove in this paper that $\nu$ is strictly more powerful than both $\ell$ and $b$, which makes $\nu$ the unique smallest reachable measure of repetitiveness to date.

There are several open questions related to NU-systems and $\nu$. For example, does it hold that $\nu = \Omega(\ell \log \log n / \log n)$, or $\nu = \Omega(\delta / \sqrt{n})$, for every string family? Is $\nu = O(\gamma)$, or at least $o(\gamma \log(n/\gamma))$, for every string family? (recall that $\gamma$ and $o(\gamma \log(n/\gamma))$ space is unknown to be reachable [9]). And towards having a practical compressor based on $\nu$, can we decompress a NU-system efficiently?

In a more general perspective, this paper pushes a little further the discussion of what we understand by a repetitive string. Intuitively, repetitiveness is about copies, and macro schemes capture those copies pretty well, but there are other aspects in a text that could be repeated besides explicit copies, such as general patterns and the relative ordering of symbols. Macro schemes capture explicit copies, L-systems capture self-similarity, and NU-systems...
capture both. What other regularities could we exploit when compressing strings, keeping the representation (more or less) simple and the associated repetitiveness measure (hopefully efficiently) computable?

References


Abstract

Suppose we are asked to index a text $T[0..n-1]$ such that, given a pattern $P[0..m-1]$, we can quickly report the maximal substrings of $P$ that each occur in $T$ at least $k$ times. We first show how we can add $O(r \log n)$ bits to Rossi et al.’s recent MONI index, where $r$ is the number of runs in the Burrows-Wheeler Transform of $T$, such that it supports such queries in $O(km \log n)$ time. We then show how, if we are given $k$ at construction time, we can reduce the query time to $O(m \log n)$.

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1 Introduction

In his foundational text Compact Data Structures: A Practical Approach [9, Section 11.6.1], Navarro posed the following problem:

“Assume we have the suffix tree of a collection of genomes $T[0..n-1]$. We then receive a short DNA sequence $P[0..m-1]$ and want to output all the maximal substrings of $P$ that appear at least $k$ times in $T$... Those substrings of $P$ are likely to have biological significance.”

He described how to solve the problem with a suffix tree for $T$ in $O(m \text{polylog}(n))$ time. Since $T$ is a collection of genomes, it is likely to be highly repetitive and the theoretically best suffix-tree implementation is likely to be the $O(r \log(n/r))$-space one by Gagie, Navarro and Prezza [5], where $r$ is the number of runs in the Burrows-Wheeler Transform (BWT) of $T$. That full data structure is complicated, however, and has never been implemented.

---

1 We have changed $T$ to $T[0..n-1]$ and $P[1..m]$ to $P[0..m-1]$ for consistency with the rest of this paper, and omitted a parameter bounding from below the length of the substrings (since we can filter them afterwards).
MONI Can Find $k$-MEMs

Very recently, Navarro [10] also gave solutions not based on a suffix tree, with the following bounds:

- $O(g_\delta)$ space and $O(km^2 \log^e n)$ query time;
- $O(\delta \log(n/\delta))$ space and $O(m \log m (\log m + k \log^e n))$ query time;
- $O(g)$ space and $O(m^2 \log^{2+\epsilon} n)$ query time when $k = \omega(\log^2 n)$;
- $O(\gamma \log(n/\gamma))$ space and $O(m \log m \log^{2+\epsilon} n)$ query time when $k = \omega(\log^2 n)$.

We refer readers to Navarro’s paper and the references therein for definitions of $g_\delta$, $\delta$, $g$ and $\gamma$.

In this paper we first show how we can add $O(r \log n)$ bits to Rossi et al.’s [13] recent MONI index to obtain a solution with $O(km \log n)$ query time. We then show how, if we are given $k$ at construction time, we can reduce the query time to $O(m \log n)$, simultaneously using less space than Gagie et al.’s compressed suffix tree and less time than Navarro’s solutions. The rest of the paper is laid out as follows: in Section 2 we review MONI in enough depth to build on it; in Section 3 we show how we can extend $\phi$ queries to support sequential access to the LCP array; in Section 4 we show how to use $\phi$ and $\phi^{-1}$ queries and LCP access to obtain a solution with $O(km \log n)$ query time; in Section 5 we show how, if we are given $k$ at construction time, we can precompute some answers, reducing the query time to $O(m \log n)$; and we conclude in Section 6. For the sake of brevity we assume readers are familiar with the concepts in Navarro’s text.

2 MONI

Bannai, Gagie and I [1] designed an index for $T$ that takes $O(r \log n)$ bits plus the space needed to support fast random access to $T$, and lists all the maximal exact matches (MEMs) of $P$ with respect to $T$—that is, all the substrings $P[i..j]$ of $P$ occurring in $T$ such that $i = 0$ or $P[i−1..j]$ does not occur in $T$, or $j = m − 1$ or $P[i..j + 1]$ does not occur in $T$—in $O(m \log \log n)$ time plus the time needed for $O(m)$ random accesses to $T$. MEMs are widely used in DNA alignment [8] and they are the substrings of $P$ Navarro asks for when $k = 1$. Generalizing to arbitrary $k$, we refer to the substrings he asks for as $k$-MEMs. Bannai et al. did not give an efficient construction algorithm or an implementation, but Rossi et al. later did. They called their implementation MONI, the Finnish word for “multi”, since it is intended to store a multi-genome reference. Boucher et al. [2] then gave a version of MONI that processes $P$ online using longest common extension (LCE) queries on $T$ instead of random access. We can support those LCE queries in $O(\log n)$ time with a balanced straight-line program for $T$, which in practice takes significantly less space than the rest of MONI.

We now sketch how Boucher et al.’s version of MONI works, incorporating ideas from Nishimoto and Tabei [12] and Brown, Gagie and Rossi [4] about replacing rank queries by table lookup and assuming we have an LCE data structure. Suppose

$$T = \text{GATTACAT#AGATACAT#GATAGAT#GATTAGA}$$

with $\$ < # < A < · · · < T, and consider Table 1, in which the permutation FL is just the inverse of the more familiar permutation LF.

For each of the value $j$ between 0 and $r − 1 = 13$, we conceptually extract from this table the starting and ending positions head$(j)$ and tail$(j)$ of run $j$ in the BWT, $\text{SA[head}(j)\text{]}$, $\text{SA[tail}(j)\text{]}$, $\text{BWT[head}(j)\text{]}$, $\text{LF[head}(j)\text{]}$ and the rank of the predecessor of $\text{LF[head}(j)\text{]}$ in the set

$$\{\text{head}[0], \ldots, \text{head}[r − 1]\}.$$
Table 1 The full table from which we conceptually start when building MONI for our example text $T = \text{GATTACAT}#AGATACAT#GATACAT#GATTAGAT#GATTAGATA}.$

<table>
<thead>
<tr>
<th>$i$</th>
<th>$\text{SA}[i]$</th>
<th>$\text{LCP}[i]$</th>
<th>lexicographically $i$th cyclic shift of $T$</th>
<th>$\text{BWT}[i]$</th>
<th>$\text{LF}(i)$</th>
<th>$\text{FL}(i)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>44</td>
<td>0</td>
<td>GATTACAT#AGATACAT#GATACAT#GATTAGAT#GATTAGATA</td>
<td>A</td>
<td>5</td>
<td>29</td>
</tr>
<tr>
<td>1</td>
<td>8</td>
<td>0</td>
<td>#GATTACAT#GATACAT#GATTAGAT#GATTAGATA#GATTACAT</td>
<td>T</td>
<td>32</td>
<td>11</td>
</tr>
<tr>
<td>2</td>
<td>17</td>
<td>1</td>
<td>#GATTACAT#GATTAGAT#GATTAGATA#GATTACAT#GATACAT</td>
<td>T</td>
<td>33</td>
<td>25</td>
</tr>
<tr>
<td>3</td>
<td>25</td>
<td>4</td>
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</table>
Table 2 The values we extract from Table 1, with the last two columns sorted.

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<th>j</th>
<th>head(j)</th>
<th>SA[head(j)]</th>
<th>tail(j)</th>
<th>SA[tail(j)]</th>
<th>BWT[head(j)]</th>
<th>μ(j)</th>
<th>finger(j)</th>
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</thead>
<tbody>
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<td>A</td>
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</tbody>
</table>

In practice we can compute the values directly without building Table 1, using prefix-free parsing [3].

We build Table 2 with these values but we sort the last two columns, which we refer to as μ(j) and finger(j). An equivalent way to define μ(j) and finger(j), illustrated in Table 1, is to draw boxes corresponding to the runs in the BWT, permute those boxes according to LF, and write their starting positions in order as the μ(j) values and the numbers of the runs in the BWT covering their starting positions as the finger(j) values. Storing Table 2 takes about

\[2r \log(n/r) + 2r \log n + r \log \sigma + 2r\]

bits, where σ is the size of the alphabet. (We do not actually need to store tail(j) = head(j + 1) − 1, of course, but we include it in Table 2 to simplify our explanation.) It is within a reasonable constant factor of the most space-efficient implementation and simple to build.

For each suffix \(P[i..m-1]\) of \(P\) from shortest to longest, MONI finds the length \(\ell_i\) of the longest prefix \(P[i..i+\ell_i-1]\) of \(P[i..m-1]\) that occurs in \(T\), the lexicographic rank \(q_i\) of a suffix of \(T\) starting with \(P[i..i+\ell_i-1]\), the starting position \(SA[q_i]\) of that suffix in \(T\), and the row \(j_i\) of Table 2 such that head\(\langle j_i \rangle\) is the predecessor of \(q_i\) in that column. We note that the (pos, len) pairs \((SA[q_0], \ell_0), \ldots, (SA[q_{m-1}], \ell_{m-1})\) are the matching statistics MS[0..m-1] of \(P\) with respect to \(T\).

Suppose we know \(i, \ell_i, q_i, SA[q_i]\) and \(j_i\), and we want to find \(\ell_{i-1}, q_{i-1}, SA[q_{i-1}]\) and \(j_{i-1}\). If BWT[head\(\langle j_i \rangle\)] = \(P[i-1]\) then we perform an LF step, as we describe in a moment. If BWT[head\(\langle j_i \rangle\)] \(\neq P[i-1]\) then we find the last row \(j_i'\) above row \(j_i\) with BWT[head\(\langle j_i' \rangle\)] = \(P[i-1]\), and the first row \(j_i''\) below row \(j_i\) with BWT[head\(\langle j_i'' \rangle\)] = \(P[i-1]\), using rank and select queries on column BWT[head\(\langle j_i \rangle\)] in Table 2. We use LCE queries to check whether \(T[SA[q_i]..n-1]\) has a longer common suffix with \(T[SA[tail(j_i')]..n-1]\) or with \(T[SA[head(j_i'')]..n-1]\) and, depending on that comparison, either reset

\[\ell_i = \text{LCE}(SA[q_i], SA[tail(j_i')])\]
\[q_i = \text{tail}(j_i')\]
\[SA[q_i] = SA[tail(j_i'')]\]
\[j_i = j_i''\]
or reset
\[
\ell_i = \text{LCE}(\text{SA}[q_i], \text{SA}[\text{head}(j_i)]) \\
q_i = \text{head}(j_i) \\
\text{SA}[q_i] = \text{SA}[\text{head}(j_i)] \\
j_i = j_i.
\]

Now \(\text{BWT}[\text{head}(j_i)] = P[i-1]\), so we can proceed with the LF step.

For example, suppose \(P[0..11] = \text{TAGATTACATTA}\), \(i = 2\) and we have already found \(\ell_2 = 8\) (because \text{GATTACAT} occurs in \(T\) but \text{GATTACATT} does not), \(q_2 = 29\), \(\text{SA}[q_2] = 0\) and \(j_2 = 7\). Since \(\text{BWT}[\text{head}(7)] = 8 \neq P[1] = A\), we find \(j_2' = 5\) and \(j_2'' = 9\) and compare
\[
\text{LCE}(0, \text{SA}[\text{tail}(5)]) = \text{LCE}(0, 10) = 3
\]
against
\[
\text{LCE}(0, \text{SA}[\text{head}(9)]) = \text{LCE}(0, 7) = 0.
\]

Since the former LCE is longer, we set \(\ell_2 = 3\), \(q_2 = 27\), \(\text{SA}[q_2] = 10\) and \(j_2 = 5\).

To perform an LF step with Table 2 when we know \(i, \ell_i, q_i, \text{SA}[q_i]\) and \(j_i\), we first set
\[
\ell_{i-1} = \ell_i + 1 \\
q_{i-1} = \mu(\pi(j_i)) + q_i - \text{head}(j_i) \\
\text{SA}[q_{i-1}] = \text{SA}[q_i] - 1,
\]
where \(\pi\) is the permutation on \(\{0, \ldots, r-1\}\) that stably sorts the column \(\text{BWT}[\text{head}(j)]\). If we keep \(\text{BWT}[\text{head}(j)]\) in a wavelet tree then we have fast access to \(\pi\).

For our example, consider \(i = 2\), \(\ell_2 = 3\), \(q_2 = 27\), \(\text{SA}[q_2] = 10\) and \(j_2 = 5\). Since \(\text{BWT}[\text{head}(5)]\) is the second \(A\) in the column \(\text{BWT}[\text{head}(j)]\) and there are 4 characters in the column lexicographically strictly less than \(A\), \(\pi(5) = 5\) and \(\mu(5) = 6\), so we set \(\ell_1 = 4\), \(q_1 = 6 + 27 - 22 = 11\) and \(\text{SA}[11] = 9\). Notice \(\pi\) is similar to an LF mapping for the sequence obtained by sampling one character from each run of the BWT (but in our example \(\pi\) has a fixed point at 5); in fact, it permutes the coloured boxes in Table 1 according to LF. It follows that \(\mu(\pi(j_i)) = \text{LF}(\text{head}(j_i))\). Since LF maintains the relationship between elements in the same box,
\[
\text{LF}(q_i) - \text{LF}(\text{head}(j_i)) = q_i - \text{head}(j_i);
\]
substituting and rearranging, we obtain our formula for \(q_{i-1}\).

The last thing left for us to do during an LF step is find \(j_{i-1}\). For this, we use the finger\((j)\) column. By construction, \(\text{head}(\text{finger}(\pi(j_i)))\) is the predecessor of \(\text{LF}(\text{head}(j_i))\) in the set
\[
\{\text{head}[0], \ldots, \text{head}[r-1]\}.
\]

Therefore, since \(q_{i-1} = \text{LF}(q_i) \geq \text{LF}(\text{head}(j_i))\), we can find the row \(j_{i-1}\) of Table 2 such that \(\text{head}(j_{i-1})\) is the predecessor of \(q_{i-1}\) in that column, by starting an exponential search at row \(\text{finger}(\pi(j_i))\). This takes \(O(\log r)\) time in the worst case and in practice it takes constant time. Nishimoto and Tabei showed how to guarantee it takes constant time at the cost of increasing the size of Table 2 slightly.

For more formal discussions, we refer readers to previous papers on MONI [13] and the \(r\)-index [5, 12, 4, 11].
Table 3 The table we use for $\phi$ queries and access to the LCP.

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<th>$j$</th>
<th>SA[head($j$)]</th>
<th>SA[tail($j$)]</th>
<th>LCP[head($j$)]</th>
<th>finger($j$)</th>
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<td>12</td>
<td>3</td>
<td>6</td>
<td>2</td>
</tr>
<tr>
<td>9</td>
<td>18</td>
<td>10</td>
<td>11</td>
<td>7</td>
</tr>
<tr>
<td>10</td>
<td>26</td>
<td>0</td>
<td>5</td>
<td>0</td>
</tr>
<tr>
<td>11</td>
<td>29</td>
<td>20</td>
<td>2</td>
<td>9</td>
</tr>
<tr>
<td>12</td>
<td>32</td>
<td>23</td>
<td>11</td>
<td>9</td>
</tr>
<tr>
<td>13</td>
<td>44</td>
<td>37</td>
<td>0</td>
<td>12</td>
</tr>
</tbody>
</table>

3 LCP access

We can support $\phi$ queries with table lookup as well: for each run BWT[$i..j$] in the BWT, we store SA[$i$] and SA[($i - 1$) mod $n$] as a row; we sort the rows by their first components; and we add to each row the number of the row containing the predecessor of the second component in the first column. Abusing notation slightly, we refer to the columns of the resulting table as SA[head($j$)], SA[tail($j$)] and finger($j$). Table 3 is for our running example, augmented with a column LCP[head($j$)] that stores the length of the longest common prefix of $T[SA[head(j)]..n - 1]$ and $T[SA[tail(j)]..n - 1]$. Since we are storing the row containing the predecessor of each entry in SA[tail($j$)] in the column SA[head($j$)], we can encode each entry in SA[tail($j$)] as the difference between it and its predecessor in SA[head($j$)].

Analysis shows the table then takes about $3r \log(n/r) + r \log r$ bits: we essentially gap-code the interleaving of column SA[head($j$)] and the sorted column SA[tail($j$)], which consists of $2r$ sorted numbers between 0 and $n - 1$ and thus takes about $2r \log(n/r)$ bits; Kärkkäinen, Kempa and Piątkowski [6] showed that the entries in LCP[head($j$)] sum to $O(n \log r)$ so, by Jensen’s Inequality, we can store them in a total of about $r \log O(n \log n) = r \log(n/r) + r \log \log r + O(r)$ bits; and finger($j$) takes about $r \log r$ bits.

To see how we use Table 3 to answer $\phi$ queries, suppose we know that the predecessor of 24 in SA[head($j$)] is in row 9. Then we have

$$\phi(24) = SA[tail(9)] + 24 - SA[head(9)] = 10 + 24 - 18 = 16.$$  

We know that the predecessor of 10 in SA[head($j$)] is in row finger(9) = 7, but the predecessor of 16 could be in a later row. Again, we perform an exponential search starting in row finger(9) = 7 and find the predecessor 12 of 16 in row 8. Then we have

$$\phi(16) = SA[tail(8)] + 16 - SA[head(8)] = 3 + 16 - 12 = 7.$$  

Looking at rows 32 to 34 in Table 1, we see that indeed $\phi(24) = 16$ and $\phi(16) = 7$. This works because, similar to the equation for LF, if BWT[$j - 1$] = BWT[$j$] then $\phi(SA[j - 1]) = \phi(SA[j]) - 1$. Again, for more formal discussions, we refer readers to previous papers on the $r$-index [5, 12, 4, 11].
We do not know how to support random access to the LCP array quickly in $O(r \log n)$ bits, but we can use Table 3 to provide a kind of sequential access to it. Specifically, as we use $\phi$ to enumerate the values in the SA – without necessarily knowing the positions of the cells of the SA those values appear in – we can use similar computations to enumerate the corresponding values in the LCP array. In our example, since the predecessor 18 of 24 in $SA[\text{head}(j)]$ is in row 9, we can compute the LCP value corresponding to the SA value 24 as

$$LCP[SA^{-1}[24]] = LCP[\text{head}(9)] + \text{SA[head}(9)] - 24 = 11 + 18 - 24 = 5.$$  

Checking this, we see that $LCP[SA^{-1}[24]] = LCP[34] = 5$. Since the predecessor 12 of $\phi(24) = 16$ in $SA[\text{head}(j)]$ is in row 8 of Table 3,

$$LCP[SA^{-1}[16]] = LCP[\text{head}(8)] + \text{SA[head}(8)] - 16 = 6 + 12 - 16 = 2.$$  

Checking this, we see that $LCP[SA^{-1}[16]] = LCP[33] = 2$.

Notice we do not use the SA row numbers 34 and 33 to compute the LCP value, as the SA value 24 is sufficient. We could avoid using the inverse suffix array $SA^{-1}$ in our formula by writing $LCP[SA^{-1}[24]]$ as $PLCP[24]$, for example, where $PLCP[0..n-1]$ denotes the permuted LCP array [7] of $T$. The kind of sequential access we obtain to the LCP is actually random access to the $PLCP$ array, and it is easier to explain why it works from that perspective – because if $BWT[j - 1] = BWT[j]$ then $PLCP[SA[j] - 1] = PLCP[SA[j]] + 1$.\footnote{The formula for PLCP has a $+1$ where the formula for $\phi$ has a $-1$,}

$$\phi^{-1}(SA[j] - 1) = \phi(SA[j]) - 1$$

$$PLCP[SA[j] - 1] = PLCP[SA[j]] + 1,$$

because if $BWT[j - 1] = BWT[j]$ then moving from $j$ to $LF(j)$ decrements the SA entry but increments the LCP entry.

Symmetric to using Table 3 to support $\phi$ queries, we can use a table to support $\phi^{-1}$ queries. In fact, the $(SA[\text{head}(j)], \text{SA[tail}(j)])$ pairs in the table are the same, but sorted by their second components; now we add to each row the number of the row containing the predecessor in the second column of the first component. Since we are storing the row containing the predecessor of each entry in $SA[\text{head}(j)]$ in the column $\text{SA[tail}(j)]$, we can encode each entry in $SA[\text{head}(j)]$ as the difference between it and its predecessor in $\text{SA[tail}(j)]$. Analysis then shows the table takes about $2r \log(n/r) + r \log r$ bits. Table 4 is for supporting $\phi^{-1}$ queries on our running example. For example, if we know that the predecessor of 7 in $\text{SA[tail}(j)]$ is in row 1, then we can compute

$$\phi^{-1}(7) = \text{SA[head}(1)] + 7 - \text{SA[tail}(1)] = 12 + 7 - 3 = 16$$

and we can find the row containing the predecessor of 16 in $\text{SA[tail}(j)]$ with an exponential search starting at row $\text{finger}(1) = 3$ (and ending in the same row). We can then compute

$$\phi^{-1}(16) = \text{SA[head}(3)] + 16 - \text{SA[tail}(3)] = 18 + 16 - 10 = 24$$

and we can find the row 6 containing the predecessor of 24 in $\text{SA[tail}(j)]$ with an exponential search starting at row $\text{finger}(3) = 4$.\footnote{The formula for PLCP has a $+1$ where the formula for $\phi$ has a $-1$,}
Table 4 The table we use for $\phi^{-1}$ queries.

<table>
<thead>
<tr>
<th>j</th>
<th>SA[head(j)]</th>
<th>SA[tail(j)]</th>
<th>finger(j)</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>26</td>
<td>0</td>
<td>6</td>
</tr>
<tr>
<td>1</td>
<td>12</td>
<td>3</td>
<td>3</td>
</tr>
<tr>
<td>2</td>
<td>6</td>
<td>9</td>
<td>1</td>
</tr>
<tr>
<td>3</td>
<td>18</td>
<td>10</td>
<td>4</td>
</tr>
<tr>
<td>4</td>
<td>0</td>
<td>18</td>
<td>0</td>
</tr>
<tr>
<td>5</td>
<td>29</td>
<td>20</td>
<td>6</td>
</tr>
<tr>
<td>6</td>
<td>32</td>
<td>23</td>
<td>6</td>
</tr>
<tr>
<td>7</td>
<td>7</td>
<td>35</td>
<td>11</td>
</tr>
<tr>
<td>8</td>
<td>5</td>
<td>36</td>
<td>1</td>
</tr>
<tr>
<td>9</td>
<td>44</td>
<td>37</td>
<td>13</td>
</tr>
<tr>
<td>10</td>
<td>2</td>
<td>38</td>
<td>0</td>
</tr>
<tr>
<td>11</td>
<td>9</td>
<td>39</td>
<td>2</td>
</tr>
<tr>
<td>12</td>
<td>3</td>
<td>42</td>
<td>1</td>
</tr>
<tr>
<td>13</td>
<td>8</td>
<td>44</td>
<td>1</td>
</tr>
</tbody>
</table>

With these two $O(r \log n)$-bit tables, given $k$, $j$ and $SA[j]$, we can compute $SA[j - k + 1..j + k - 1]$ and $LCP[j - k + 1..j + k - 1]$ in $O(k \log r) \subseteq O(k \log n)$ time. (Actually, we can achieve that bound even without the $\phi(j)$ columns in the tables, but Brown et al.’s results suggest those will provide a significant speedup in practice.) With Nishimoto and Tabei’s modification, we can reduce that to $O(k)$ time while keeping the tables in $O(r \log n)$ bits; this would slightly improve the time bound we give in the next section to $O(m(k + \log n))$.

Lemma 1. We can store two $O(r \log n)$-bit tables such that, given $k$, $j$ and $SA[j]$, we can compute $SA[j - k + 1..j + k - 1]$ and $LCP[j - k + 1..j + k - 1]$ in $O(k \log n)$ time.

4 Finding $k$-MEMs with Lemma 1

We store the tables described in Sections 2 and 3 for $T$, which add $O(r \log n)$ bits to MONI. Given $P$ and $k$, we find the MEMs of $P$ with respect to $T$ as before but then, from each $SA[q_i]$, we use Lemma 1 to find $LCP[q_i - k + 2..q_i + k - 1]$ in $O(k \log n)$ time.

For example, suppose that $P[0..11] = \text{TAGATTACAT}$, as in Section 2, and $k = 3$. Starting with $q_{12} = 22$, with MONI we compute the values shown in columns $q_i$, $SA[q_i]$, $\ell_i$ and $BWT[q_i]$ of Table 5. (It is important that we choose $q_i$ to be one of the endpoints of a run, since we store SA entries only at those positions, but this is true also for MONI.) The crossed out values are the ones we replace because $BWT[q_i] \neq P[i]$. If we look at the original SA$[q_i]$ and $\ell_i$ values, before any replacements, we obtain the matching statistics

\[
\text{MS}[0..11] = (38, 5), (9, 4), (0, 8), (1, 7), (2, 6), (20, 5), (21, 4), (22, 3), (1, 4), (2, 3), (3, 2), (4, 1)
\]

of $P$ with respect to $T$, with $(\text{pos, len})$ pair $\text{MS}[i]$ indicating the starting position $\text{MS}[i]$.pos in $T$ of an occurrence of the longest prefix of $P[i..m - 1]$ that occurs in $T$, and the length $\text{MS}[i]$.len of that prefix.

From the matching statistics, it is easy to compute the MEMs $P[0..4] = \text{TAGAT}$, $P[2..9] = \text{GATTACAT}$ and $P[8..11] = \text{ATTA}$ of $P$ with respect to $T$: a MEM starts at any position $i$ such that $i = 0$ or $\text{MS}[i - 1]$.len $\leq \text{MS}[i]$.len. For each $i$, after we compute $q_i$, $SA[q_i]$ and $\ell_i$ (and before we replace them, if we do), we use Lemma 1 to compute the sub-interval $LCP[q_i - 1..q_i + 2]$ of length $4 = 2k - 2$. 
Table 5 With MONI we compute the values shown in columns \( q_i \), \( \text{SA}[q_i] \), \( \ell_i \), and \( \text{BWT}[q_i] \) on the left side of the table, and from those we can compute the matching statistics and MEMs of \( P[0..11] = \text{TAGATTACATTA} \) with respect to \( T[0..44] = \text{GATTACAT#AGATACAT#GATACAT#GATTAGAT#GATTAGAT#} \). After we have computed the values on the left side of the table, we can also compute the values in columns \( \text{LCP}[q_i - k + 2..q_i + k - 1] \), \( L_i \), and \( \min(\ell_i, L_i) \) on the right side of the table, and from those we can compute the 3-MEMs of \( P \) with respect to \( T \).

<table>
<thead>
<tr>
<th>( i )</th>
<th>( P[i] )</th>
<th>( q_i )</th>
<th>( \text{SA}[q_i] )</th>
<th>( \ell_i )</th>
<th>( \text{BWT}[q_i] )</th>
<th>( \text{LCP}[q_i - 1..q_i + 2] )</th>
<th>( L_i )</th>
<th>( \min(\ell_i, L_i) )</th>
</tr>
</thead>
<tbody>
<tr>
<td>12</td>
<td>22</td>
<td>5</td>
<td>0</td>
<td>A</td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>11</td>
<td>A</td>
<td>6</td>
<td>4</td>
<td>1</td>
<td>T</td>
<td>[0, 1, 5, 8]</td>
<td>5</td>
<td>1</td>
</tr>
<tr>
<td>10</td>
<td>T</td>
<td>37</td>
<td>3</td>
<td>2</td>
<td>T</td>
<td>[1, 2, 6, 9]</td>
<td>6</td>
<td>2</td>
</tr>
<tr>
<td>9</td>
<td>T</td>
<td>42</td>
<td>2</td>
<td>3</td>
<td>A</td>
<td>[5, 1, 3, 6]</td>
<td>3</td>
<td>3</td>
</tr>
<tr>
<td>8</td>
<td>A</td>
<td>14</td>
<td>12</td>
<td>2</td>
<td>C</td>
<td>[0, 2, 4, 7]</td>
<td>4</td>
<td>4</td>
</tr>
<tr>
<td>7</td>
<td>C</td>
<td>24</td>
<td>22</td>
<td>3</td>
<td>A</td>
<td>[4, 7, 0, 3]</td>
<td>4</td>
<td>3</td>
</tr>
<tr>
<td>6</td>
<td>A</td>
<td>8</td>
<td>21</td>
<td>4</td>
<td>T</td>
<td>[5, 8, 1, 4]</td>
<td>5</td>
<td>4</td>
</tr>
<tr>
<td>5</td>
<td>T</td>
<td>37</td>
<td>23</td>
<td>5</td>
<td>T</td>
<td>[6, 9, 2, 5]</td>
<td>6</td>
<td>5</td>
</tr>
<tr>
<td>4</td>
<td>T</td>
<td>42</td>
<td>2</td>
<td>6</td>
<td>A</td>
<td>[5, 1, 3, 6]</td>
<td>3</td>
<td>3</td>
</tr>
<tr>
<td>3</td>
<td>A</td>
<td>19</td>
<td>1</td>
<td>7</td>
<td>G</td>
<td>[10, 2, 4, 7]</td>
<td>4</td>
<td>4</td>
</tr>
<tr>
<td>1</td>
<td>A</td>
<td>10</td>
<td>39</td>
<td>4</td>
<td>T</td>
<td>[4, 5, 1, 3]</td>
<td>4</td>
<td>4</td>
</tr>
<tr>
<td>0</td>
<td>T</td>
<td>41</td>
<td>38</td>
<td>5</td>
<td>T</td>
<td>[2, 5, 1, 3]</td>
<td>2</td>
<td>2</td>
</tr>
</tbody>
</table>

We scan each interval \( \text{LCP}[q_i - k + 2..q_i + k - 1] \) in \( O(k) \) time and find a sub-interval of length \( k - 1 \) such that the minimum LCP value \( L_i \) in that sub-interval is maximized. This LCP sub-interval corresponds to a sub-interval of length \( k \) in \( \text{SA}[q_i - k + 1..q_i + k - 1] \) containing the starting positions of \( k \) suffixes of \( T \) including \( T[\text{SA}[q_i]..n - 1] \) itself whose common prefix with \( T[\text{SA}[q_i]..n - 1] \) has the maximum possible length \( L_i \).

In our example, we can scan each interval in column \( \text{LCP}[q_i - 1..q_i + 2] \) of Table 5 and find the sub-interval of length 2 such that the minimum LCP value \( L_i \) is maximized. If we check Table 1, we find that the longest prefix of \( T[4..44] = \text{ACAT#AGATA...} \) that occurs at least 3 times in \( T \) indeed has length \( L_{14} = 5 \), the longest prefix of \( T[3..44] = \text{TACAT#AGATA...} \) that occurs at least 3 times in \( T \) indeed has length \( L_{10} = 6 \), and so on.

Since the common prefix of \( P[i..m - 1] \) and \( T[\text{SA}[q_i]] \) has the maximum possible length \( \ell_i \), the longest prefix of \( P[i..m - 1] \) that occurs at least \( k \) times in \( T \) has length \( \min(\ell_i, L_i) \). Computing \( \min(\ell_i, L_i) \) for each \( i \) takes a total of \( O(km \log n) \) time. The values \( \min(\ell_i, L_i) \) are something like a parameterized version of the lengths in the matching statistics: \( \min(\ell_i, L_i) \) is the length of the longest prefix of \( P[i..m - 1] \) that occurs at least \( k \) times in \( T \).

We can compute the \( k \)-MEMs of \( P \) with respect to \( T \) from the \( \min(\ell_i, L_i) \) values in the same way we compute MEMs from the lengths in the matching statistics: a \( k \)-MEM starts at any position \( i \) such that \( i = 0 \) or \( \min(\ell_{i-1}, L_{i-1}) \leq \min(\ell_i, L_i) \). In our example,

\[
\begin{align*}
\min(\ell_0, \ell_0') & \leq \min(\ell_1, \ell_1') = 4 \\
\min(\ell_1, \ell_1') & \leq \min(\ell_2, \ell_2') = 5 \\
\min(\ell_4, \ell_4') & \leq \min(\ell_5, \ell_5') = 5 \\
\min(\ell_7, \ell_7') & \leq \min(\ell_8, \ell_8') = 4
\end{align*}
\]

and so the \( k \)-MEMs are \( P[0..1] = \text{TA} \), \( P[1..4] = \text{AGAT} \), \( P[2..6] = \text{GATTA} \), \( P[5..9] = \text{TACAT} \) and \( P[8..11] = \text{ATTA} \).

We can compute \( \min(\ell_i, L_i) \) as soon as we have computed \( \text{SA}[q_i] \) and \( \ell_i \), so we can compute the \( k \)-MEMs of \( P \) with respect to \( T \) online.
Theorem 2. Suppose we have MONI for a text $T[0..n-1]$ whose BWT consists of $r$ runs. We can add $O(r \log n)$ bits to MONI such that, given $P[0..m-1]$ and $k$, we can find the $k$-MEMs of $P$ with respect to $T$ online in $O(k \log n)$ time per character of $P$.

Finding $k$-MEMs with precomputed values

Suppose the interval of length $k$ that we find in SA for $P[i..m-1]$, following the procedures in Section 4, is $SA[s_i..s_i + k - 1]$ and BWT[$s_i] = \cdots = BWT[s_i + k - 1] = P[i - 1]$. Then \(\min(\ell_{i-1}, L_{i-1}) = \min(\ell_i, L_i) + 1\) and we can find the interval for $P[i - 1..m - 1]$ with an LF query for $s_i$, in $O(\log n)$ time. This means we need the results of Section 3 only when at least one character in BWT[$s_i..s_i + k - 1$] is not equal to $P[i - 1]$.

First, suppose BWT[$q_i] \neq P[i - 1]$. Following the procedures in Section 2, MONI resets $q_i$ to the endpoint $b$ of a run in the BWT, resets $\ell_i$, and then computes $q_{i-1} = LF(b)$. Following the procedures in Section 4, we compute LCP[$q_{i-1} - k + 2..q_{i-1} + k - 1$] and scan it to compute the interval $SA[s_{i-1}..s_{i-1} + k - 1]$ for $P[i - 1..m - 1]$. If we are given $k$ at construction time, however, then for every endpoint $b$ of a run in the BWT, we can precompute

- the sub-interval of length $k - 1$ of $LCP[LF(b) - k + 2..LF(b) + k - 1]$ that maximizes the minimum value $L(b)$ in the sub-interval,
- that value $L(b)$.

With this information, we do not need the results of Section 3 for this case either, and can handle it in $O(\log n)$ time as well. Since the sub-interval we store for $b$ starts between $LF(b) - k + 2$ and $LF(b) + k - 1$, we can store it in $O(\log k)$ bits as an offset. This means we store $O(r \log k)$ bits on top of at most $2r$ LCP values, or $O(r \log n)$ bits in total.

The remaining case is when BWT[$q_i] = P[i - 1]$ but some of the other characters in BWT[$s_i..s_i + k - 1$] are not equal to $P[i - 1]$. If BWT[$q_i$] is the end of a run, then we can proceed as in the previous case in $O(\log n)$ time, using our precomputed values for $q_i$ (but without resetting $q_i$ and $\ell_i$). Otherwise, we claim we can choose such a character BWT[$b] = P[i - 1]$ at the end of a run, set

$$\ell_i = \min(LCE(SA[q_i], SA(b)), \ell_i)$$

and $q_i = b$, and then proceed as in the previous case in $O(\log n)$ time, and still be sure of obtaining the correct $k$-MEMs of $P$ with respect to $T$. (Continuing to run MONI with the new values of $q_i$ and $\ell_i$ may not give us the correct MEMs, however.) To be able to change $q_i$ and $\ell_i$ this way, it is important that we now work online, instead of running MONI on $P$ and then using the results to find the $k$-MEMs.

To see why our claim holds, assume our query has worked correctly so far, so

$$T[SA[q_i]..SA[q_i] + \min(\ell_i, L_i) - 1] = T[SA[b]..SA[b] + \min(\ell_i, L_i) - 1]$$

is the longest prefix of $P[i..m - 1]$ that occurs at least $k$ times in $T$. Therefore, the $k$-MEMs starting in $P[0..i - 1]$ are all completely contained in $P[0..i + \min(\ell_i, L_i) - 1]$. It follows that resetting

$$\ell_i = \min(LCE(SA[q_i], SA(b)), \ell_i)$$

and $q_i = b$ does not affect the set of $k$-MEMs we find that start in $P[0..i - 1]$.
if \(\text{BWT}[s_i] = \cdots = \text{BWT}[s_i + k - 1] = P[i - 1]\) then
\[
q_{i-1} \leftarrow \text{LF}(q_i) \\
\ell_{i-1} \leftarrow \ell_i + 1 \\
L_{i-1} \leftarrow L_i + 1 \\
s_{i-1} = \text{LF}(s_i)
\]
else
if \(\text{BWT}[q_i] \neq P[i - 1]\) then
reset \(q_i\) and \(\ell_i\) as MONI does
else if \(\text{BWT}[q_i]\) is not at the end of a run
choose \(b\) in \([s_i..s_i + k - 1]\) with \(\text{BWT}[b] = P[i - 1]\) at the end of a run
\[
\ell_i \leftarrow \min(\text{LCE}(\text{SA}[q_i], \text{SA}[b]), \ell_i) \\
q_i \leftarrow b
\]
end if
\[
q_{i-1} = \text{LF}(q_i) \\
\ell_{i-1} \leftarrow \ell_i + 1 \\
L_{i-1} \leftarrow L(q_i) \\
s_{i-1} = \text{LF}(q_i) - \text{offset}(q_i)
\]
end if

Figure 1 Pseudo-code for how we find \(k\)-MEMs with precomputed values.

Figure 1 shows pseudo-code for how we find \(k\)-MEMs with precomputed values. Table 6 shows the offsets and \(L(b)\) values for our example, surrounded by coloured boxes on the right, with each offset indicating how far above \(\text{LF}(b)\) the sub-interval starts. The coloured boxes on the left indicate the sub-interval itself and the longest common prefix of the suffixes starting in the sub-interval of the SA.

For our example, suppose we again start with \(q_{12} = 22\) and \(\ell_{12} = 0\). Since \(\text{BWT}[q_{12}] = P[11] = A\), we set \(q_{11} = \text{LF}(22) = 6\) and \(\ell_{11} = \ell_{12} + 1 = 1\). The values \(\text{offset}(22) = 0\) and \(L(22) = 5\) in the black rectangle in Table 6 tell us to set \(s_{11} = \text{LF}(22) - 0 = 6\) and \(L_{11} = 5\). This means the suffixes of \(T\) with starting points in
\[
\text{SA}[6..8] = [4, 13, 21]
\]
have a longest common prefix of length 5, which starts with the longest prefix of \(P[11]\) that occurs at least 3 times in \(T\). This longest prefix has length \(\min(\ell_{11}, L_{11}) = 1\) so it is just \(P[11] = A\). After this initial setup, we can fill in Table 7 according to the pseudo-code in Figure 1, with crossed out values again indicating those that are replaced.

\textbf{Theorem 3.} Suppose we have MONI for a text \(T[0..n - 1]\) whose BWT consists of \(r\) runs. Given \(k\), we can add \(O(r \log n)\) bits to MONI such that, given \(P[0..m - 1]\), we can find the \(k\)-MEMs of \(P\) with respect to \(T\) online in \(O(k \log n)\) time per character of \(P\).

\section{Conclusion}

We have shown, first, how we can add \(O(r \log n)\) bits to MONI for a text \(T[0..n - 1]\), where \(r\) is the number of runs in the BWT of \(T\), such that if we are given \(k\) at query time with \(P[0..m - 1]\), then we can find the \(k\)-MEMs of \(P\) with respect to \(T\) online in \(O(k \log n)\) time per character of \(P\). We have then shown how, if we are given \(k\) at construction time, we can add \(O(r \log k)\) bits and at most \(2r\) LCP values – which are \(O(r \log n)\) bits in total – such
The table showing the precomputed values we use to find 3-MEMs with respect to our example $T = GATTACAT\#AGATACAT\#GATACAT\#GATTAGAT\#GATTAGATA$.

<table>
<thead>
<tr>
<th>$i$</th>
<th>SA[$i$]</th>
<th>LCP[$i$]</th>
<th>lexicographically $i$th cyclic shift of $T$</th>
<th>BWT[$i$]</th>
<th>LF($i$)</th>
<th>offset($i$)</th>
<th>$L(i)$</th>
</tr>
</thead>
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<td>0</td>
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</tr>
</tbody>
</table>
Table 7 The values we compute (except \(\text{BWT}[s_i..s_i + 2]\), which we include here only for clarity) while finding the 3-MEMs of \(P[0..11] = \text{TAGATTACATTA}\) with respect to our example \(T = \text{GATTACAT#AGATACAT#GATACAT#GATTAGAT#GATTAGAT}\$.

<table>
<thead>
<tr>
<th>(i)</th>
<th>(q_i)</th>
<th>(\ell_i)</th>
<th>(L_i)</th>
<th>(\min(\ell_i, L_i))</th>
<th>(s_i)</th>
<th>(P[i-1])</th>
<th>(\text{BWT}[q_i])</th>
<th>(\text{BWT}[s_i..s_i + 2])</th>
</tr>
</thead>
<tbody>
<tr>
<td>12</td>
<td>22</td>
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<td></td>
<td>A</td>
<td>A</td>
<td></td>
<td></td>
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</tr>
<tr>
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<td>1</td>
<td>6</td>
<td>T</td>
<td>T</td>
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</tr>
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<td>37</td>
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<td>6</td>
<td>2</td>
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<td>3</td>
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</table>

that we can find the \(k\)-MEMs of \(P\) with respect to \(T\) online in \(O(\log n)\) time per character of \(P\). Along the way, we have also shown how to extend \(\phi\) queries to support sequential access to the LCP, which may be of independent interest.

Although we have not discussed construction, we expect it will not be difficult to modify prefix-free parsing [2] to build our tables for \(\phi\), LCP and \(\phi^{-1}\) queries. Once we can support those queries, we can use them to compute \(k\)-MEMs in \(O(km \log n)\) time, or to build in \(O(kr)\) time the table of precomputed values that we need to compute \(k\)-MEMs in \(O(m \log n)\) time. In fact, once we have built the tables for \(\phi\), LCP and \(\phi^{-1}\) queries – which take \(O(r)\) space but may be significantly larger than our table of precomputed values – then we can store them in external memory and recover them only when we want to build a table of precomputed values for a different choice of \(k\).

We believe our approach is a practical extension of MONI and we are currently implementing it. One possible application might be to index two genomic databases (possibly with two different values of \(k\)), one of haplotypes from people with symptoms of a genetic disease and one of haplotypes from people without; then, as the first step in a bioinformatics pipeline, we could use those indexes to mine for substrings that are common in one database and not in the other. We think this application is interesting because, except for a remark in Bannai et al.’s paper about potentially applying MEM-finding to rare-disease diagnosis, the \(r\)-index and MONI have so far been considered only as tools for pangenomic alignment, and this is an application to pangenomic analysis. If the disease is recessive or multifactorial then variations associated with it are likely to be present in both databases, so MEM-finding is unlikely to detect them; those variations could be more frequent in the first database, however, so \(k\)-MEM-finding may still be useful.

References


